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# **Contamination Assessment Report**

for **Site 333** 



Panama City, Florida



# Southern Division Naval Facilities Engineering Command

Contract Number N62467-94-D-0888
Contract Task Order 0008

February 1997

# CONTAMINATION ASSESSMENT REPORT FOR SITE 333

# COASTAL SYSTEMS STATION PANAMA CITY, FLORIDA

# COMPREHENSIVE LONG-TERM ENVIRONMENTAL ACTION NAVY (CLEAN) CONTRACT

Submitted to:
Southern Division
Naval Facilities Engineering Command
2155 Eagle Drive
North Charleston, South Carolina 29406

Submitted by:
Brown & Root Environmental
661 Andersen Drive
Foster Plaza 7
Pittsburgh, Pennsylvania 15220

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PREPARED BY:

APPROVED FOR SUBMITTAL BY:

GERALD F. GOODE, P.G.
TASK ORDER MANAGER
FLORIDA LICENSE NO. 0001276
BROWN & ROOT ENVIRONMENTAL
TALLAHASSEE, FLORIDA

PROGRAM MANAGER
BROWN & ROOT ENVIRONMENTAL

PITTSBURGH, PENNSYLVANIA

#### **EXECUTIVE SUMMARY**

Brown & Root Environmental (B&R Environmental) has completed a Contamination Assessment (CA) at the Coastal Systems Station (CSS) Site 333 in accordance with the requirements of Chapter 62-770, Florida Administrative Code (FAC). The Contamination Assessment Report for Site 333 completed by B & R Environmental, January 1997, was submitted to the Florida Department of Environmental Protection (FDEP) for approval.

B&R Environmental performed the following tasks during the CA:

- Reviewed available Navy documents to identify potential sources and receptors for petroleum hydrocarbons in the vicinity, to evaluate private potable wells in a 0.25-mile radius and public supply water supply wells within 0.50-mile radius, and to locate nearby surface water bodies and determine surface hydrology;
- Reviewed soil and groundwater data collected during removal of the oil/water separator and waste oil tank to determine boring locations and monitoring well placements;
- Conducted site survey to identify utilities and to construct a site plan;
- Performed excavation of 16 soil borings for organic vapor analysis;
- Installed five shallow temporary groundwater monitoring points and collected groundwater samples for field screening with a portable gas chromatograph;
- Advanced four shallow permanent monitoring wells to approximately 13 feet below land surface (bls);
- Collected groundwater samples from the permanent monitoring wells for laboratory analysis for Gasoline and Kerosene Analytical Group parameters, including used oil parameter testing;

The results of the CA identified no "excessively contaminated" soil at the site, as defined by Chapter 62-770.200 FAC. Groundwater concentrations of total naphthalene, lead, vinyl chloride, phenanthrene, cis 1-2 dichlorothene, and Total Recoverable Petroleum Hydrocarbons (TRPH) were reported at concentrations above State Action Levels, however, these constituents were at levels which meet the criteria for Monitoring Only status. In an effort to further reduce these concentration levels to meet No Further Action status, the development of an Alternative Remedial Procedure for review and approval by the FDEP is recommended prior to developing and/or implementing a Monitoring Only Plan.

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#### 1.0 INTRODUCTION

#### 1.1 PURPOSE AND SCOPE

A Contamination Assessment (CA) has been conducted by Brown and Root Environmental (B&R Environmental) for the U.S. Navy (Navy) Southern Division Naval Facilities Engineering Command under Contract Task Order 0008, for the Comprehensive Long-term Environmental Action Navy (CLEAN III), Contract Number N62467-94-D-0888. The CA was conducted at Site 333 located at the Coastal Systems Station (CSS) in Panama City, Florida. The Florida Department of Environmental Protection (FDEP) Facility Identification Number is 038518667.

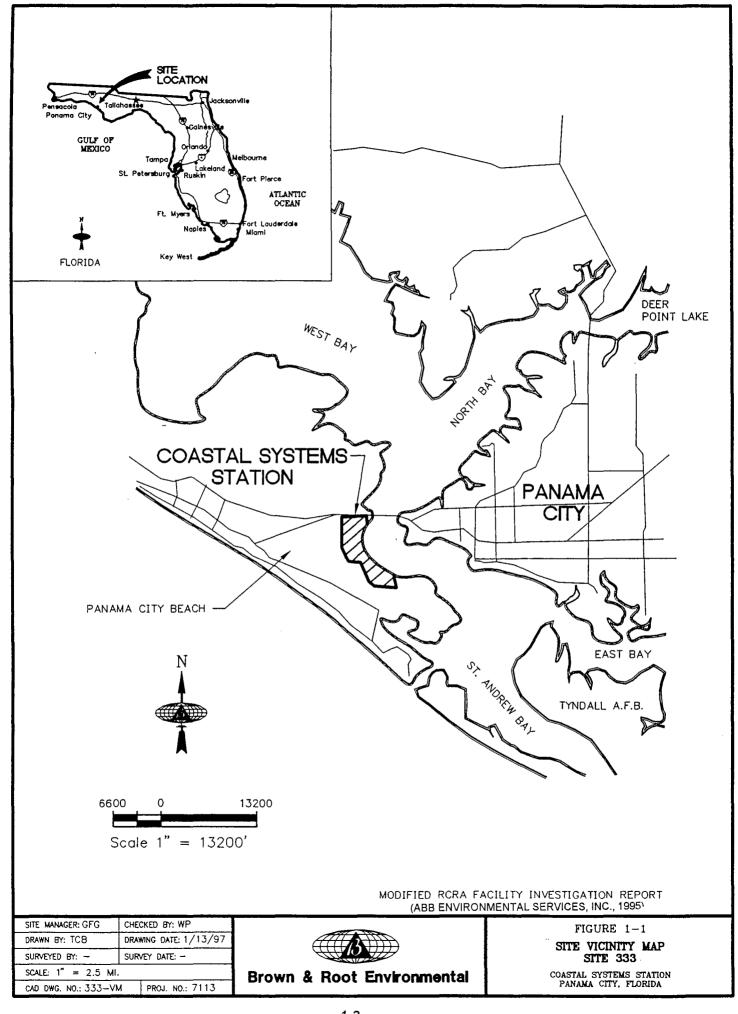
The purpose of this CA was to determine the nature and extent of petroleum hydrocarbons and used oil constituents in soil and groundwater in accordance with the requirements of Chapter 62-770 of the Florida Administrative Code (FAC). The Navy submitted a Discharge Notification Form (DNF) to the Bay County Health and Rehabilitative Services (HRS), Environmental Health Services, Pollution Storage Tank Program in December 1995. The discharge was reported based on soil vapor readings and groundwater quality testing conducted during removal of the site's waste oil tank and oil/water separator. The DNF listed the type of substance used as waste oil. The cause of the leak was reported on the DNF as unknown. A copy of the DNF report is included as Appendix A.

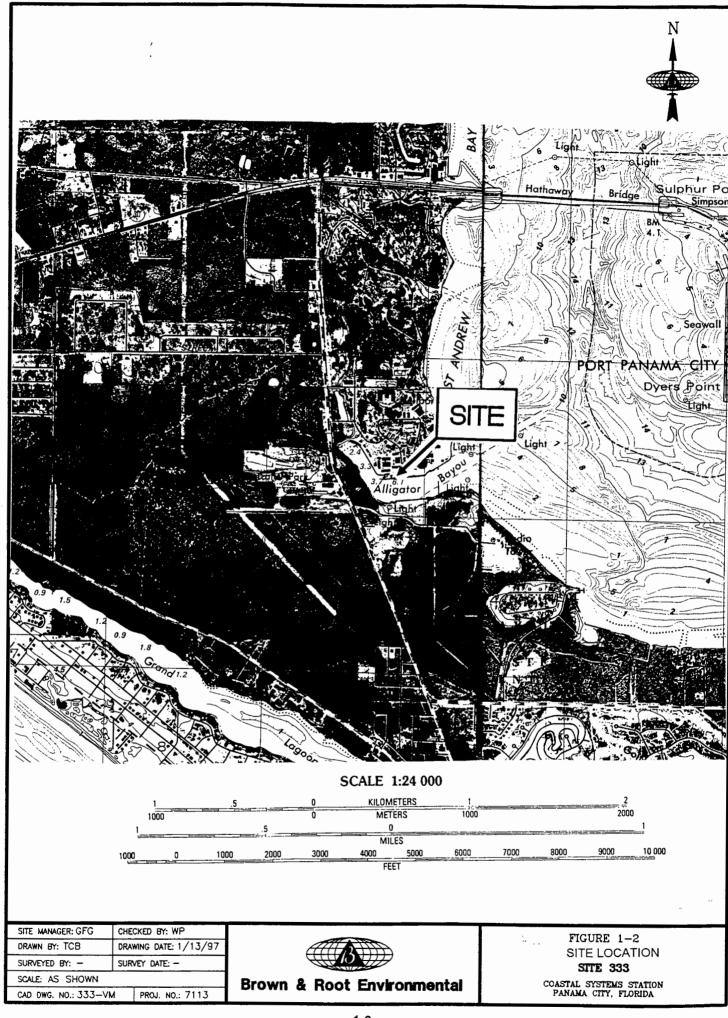
A CAR Summary Sheet, as required by Chapter 62-770, FAC., is included as Appendix B.

#### 1.2 SITE DESCRIPTION

#### 1.2.1 Location

The CSS facility is located on the western shore of St. Andrew Bay in Panama City, Bay County, Florida. The facility is bounded by U.S. Highway 98 to the north, St. Andrews Bay to the east, State Road 292B (Magnolia Beach Road) to the south and State Road 292 (Thomas Drive) to the west as shown on Figure 1-1. Specifically, the CSS facility is located within Section 33 of Township 3 South, Range 15 West and Section 4 of Township 4 South, Range 15 West, as shown on the United States Geological Survey (USGS) 7.5 Series (Topographic) Panama City Beach Quadrangle, Florida, presented as Figure 1-2.





#### 1.2.2 Topography and Surface Water

The topography at the site is relatively flat with a gentle surface slope towards the south. The site is located at an elevation of approximately 7 feet above mean sea level. The nearest surface water body is Alligator Bayou located approximately 75 feet south of the site. Alligator Bayou is designated as a Class III surface water by the State of Florida, suitable for fish and wildlife propagation and water sports (ABB Environmental Services Inc., RCRA Facility Investigation, 1995).

#### 1.2.3 Regional Hydrogeology

The regional hydrogeology of CSS Panama City is described in the RCRA Facility Investigation report (ABB Environmental Services, Inc., 1995). According to this report, surficial deposits at CSS are Pleistocene to Recent coastal plain sediment of marine and estuarine origin. They predominately consist of quartz sand, clayey sand, and gravel. These deposits vary in thickness from 70 to 100 feet in Bay County. The surficial aquifer is located within these deposits.

Underlying the surficial deposits is the Intercoastal Formation of middle Miocene to late Pliocene. The Intercoastal Formation is composed of sand and poorly consolidated limestone interbedded with discontinuous clay and low permeability sandy limestone. This formation is approximately 150 feet thick at CSS Panama City. The lower beds of the Intercoastal Formation are part of the Floridan aquifer system.

Groundwater at CSS occurs in two major aquifer systems: unconfined surficial aquifer and the Floridan aquifer system, which is under confined and artesian conditions. A third semi confined aquifer exists in thin permeable sand and shell zones within the Intercoastal Formation, and is separated from the water table aquifer and from the Floridan aquifer system by interbedded low-permeability clay and limestone. The Intercoastal Formation does not produce enough water to be considered a significant water source. The Floridan aquifer is under confined and artesian conditions where low-permeable clays and limestone beds of the Intracoastal Formation separate the water table aquifer from the Floridan aquifer. The surficial aquifer is reported to have insufficient thickness to produce significant quantities of water and its quality is generally undesirable for human use (i.e., dissolved solids, acidity, and iron content). Low permeability clay lenses in the surficial aquifer and the Intercoastal Formation are discontinuous. The surficial aquifer may be hydraulically connected to the Floridan aquifer system through semiconfining strata of the Intercoastal Formation.

#### **1.2.4 Land Use**

Site 333 is located in the southwest area of the CSS property as shown on Figure 1-3. This area of the Base is comprised of research facilities and various support activities. Potential sources of contamination near the site include a 6,000 gallon diesel underground storage tank (UST), CSS Tank No. 307R, hazardous drum storage structure 307, and Solid Waste Management Unit No. 1 (SWMU-1).

The 6,000 gallon double-walled fiberglass diesel fuel UST (CSS Tank No. 307R) includes double-walled product piping and is equipped with leak detection sensors. This tank provides diesel fuel to a dispenser located east of the boom containment wash area. The 6,000 gallon diesel UST was installed during the late 1980's as a system upgrade to replace a 5,000 gallon diesel UST. The 6,000 gallon diesel UST is located adjacent to the southeast end of the containment boom wash area. The location of 6,000 gallon diesel UST is shown on Figure 1-4.

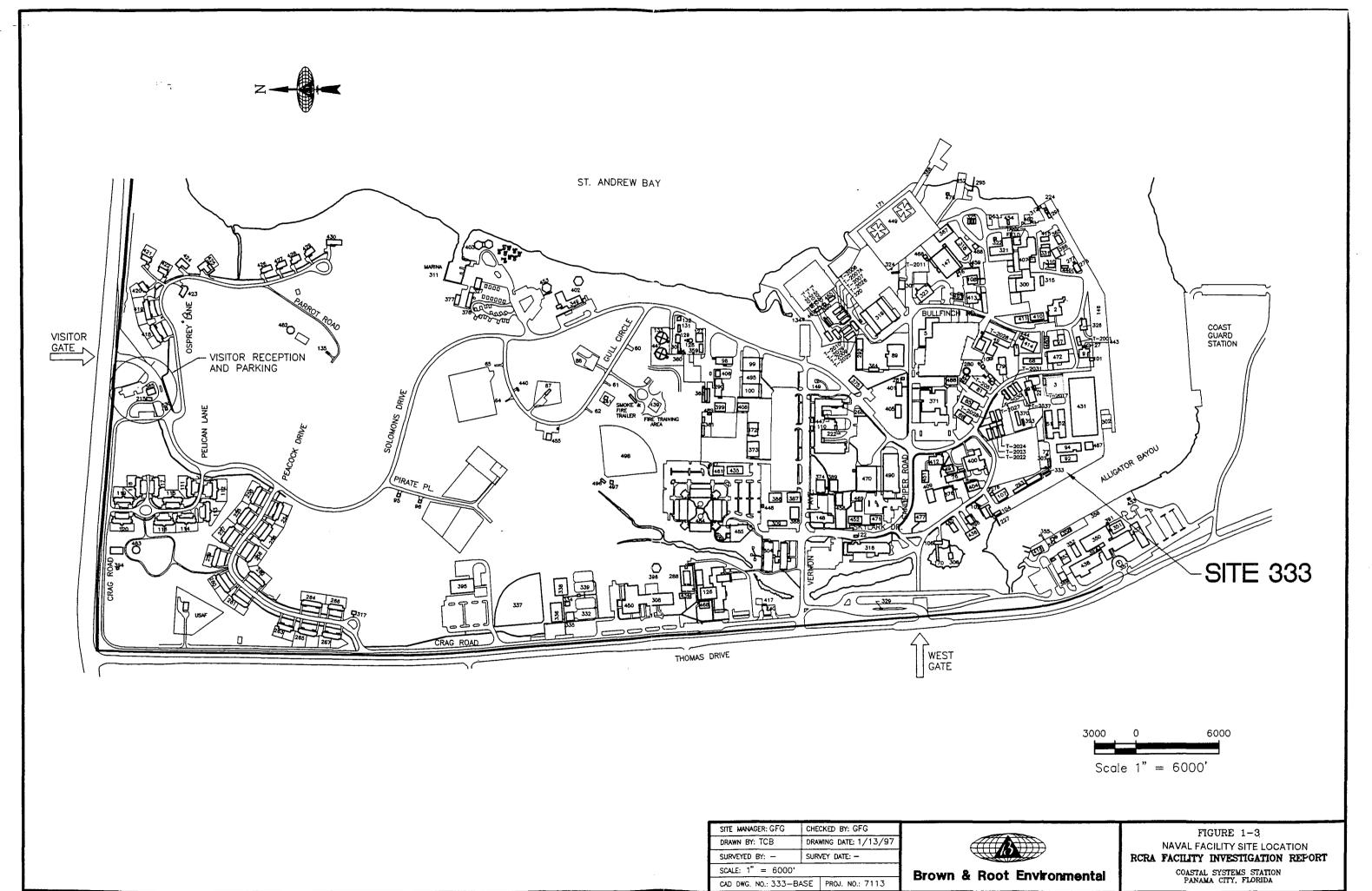
A hazardous drum storage area is contained in Structure 307. Structure 307 is secured with a locking chain link fence to restrict access to the area. The top of the structure contains a roof to shelter the drums from the sun and rain. The bottom of the structure contains concrete flooring with a floor drain to control and collect any spills which may occur. Structure 307 is located southeast of the boom containment wash area as shown on Figure 1-4.

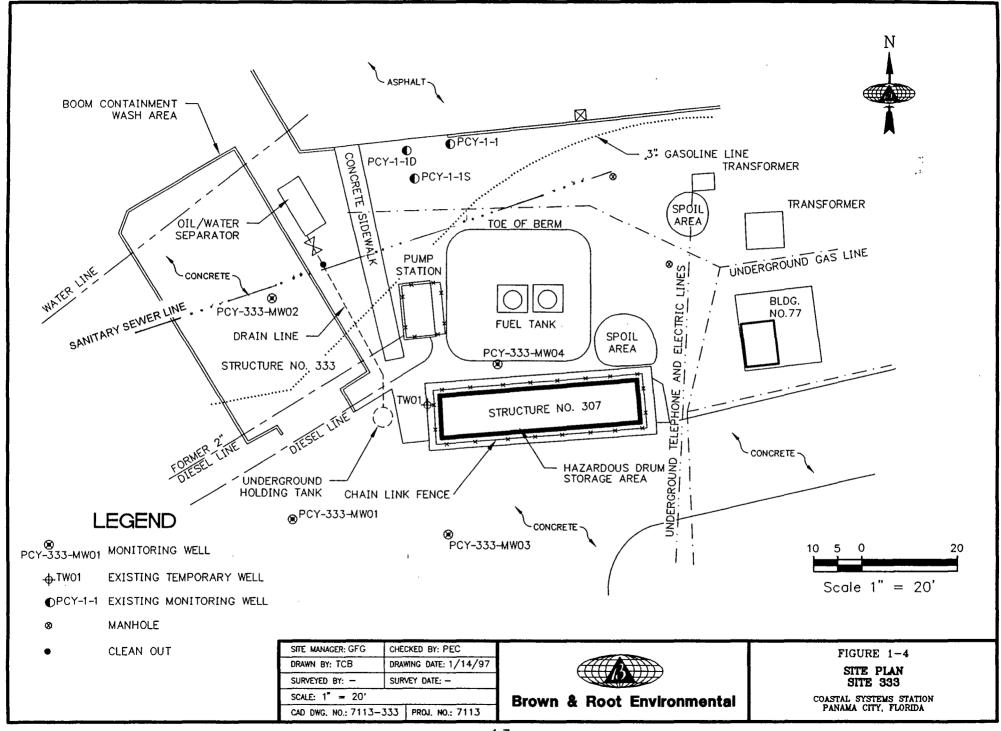
The SWMU-1 is reported to extend beneath the boom containment wash area. SWMU-1 was a marshy depression which was used as a disposal area from approximately 1945 to 1953. This disposal area received general house hold wastes, food scraps, scrap metal, scrap lumber, and small quantities of paint, paint thinner, battery acids, solvents, and photographic chemicals. Waste oil and bilge water were also poured on the ground and burned (ABB Environmental Services Inc., 1995).

Underground utilities and petroleum distribution lines identified near Site 333 include diesel lines from the adjacent diesel UST, a 3-inch diameter gasoline line, sanitary sewer, gas, telephone, and water. Underground utilities are shown on Figure 1-4.

#### 1.2.5 Site Description

Site 333 is a boom containment wash area. The site is primarily used as a containment area for the cleaning of booms used in the containment of diesel spills. An above ground 100 gallon per minute capacity oil/water separator and an underground 550-gallon fiberglass waste oil tank (underground





holding tank) are used to collect and process rinse water from the cleaning of the booms. Rinse water from the containment area is drained into sumps then pumped to the oil/water separator. Any oils collected in the separator are gravity drained into the waste oil tank. Water in the oil/water separator then drains into the Navy's sanitary sewer system. The oil/water separator also receive water from a floor drain in structure 307. Existing site features are shown on Figure 1-4.

#### 1.2.6 Potable Water Well Survey

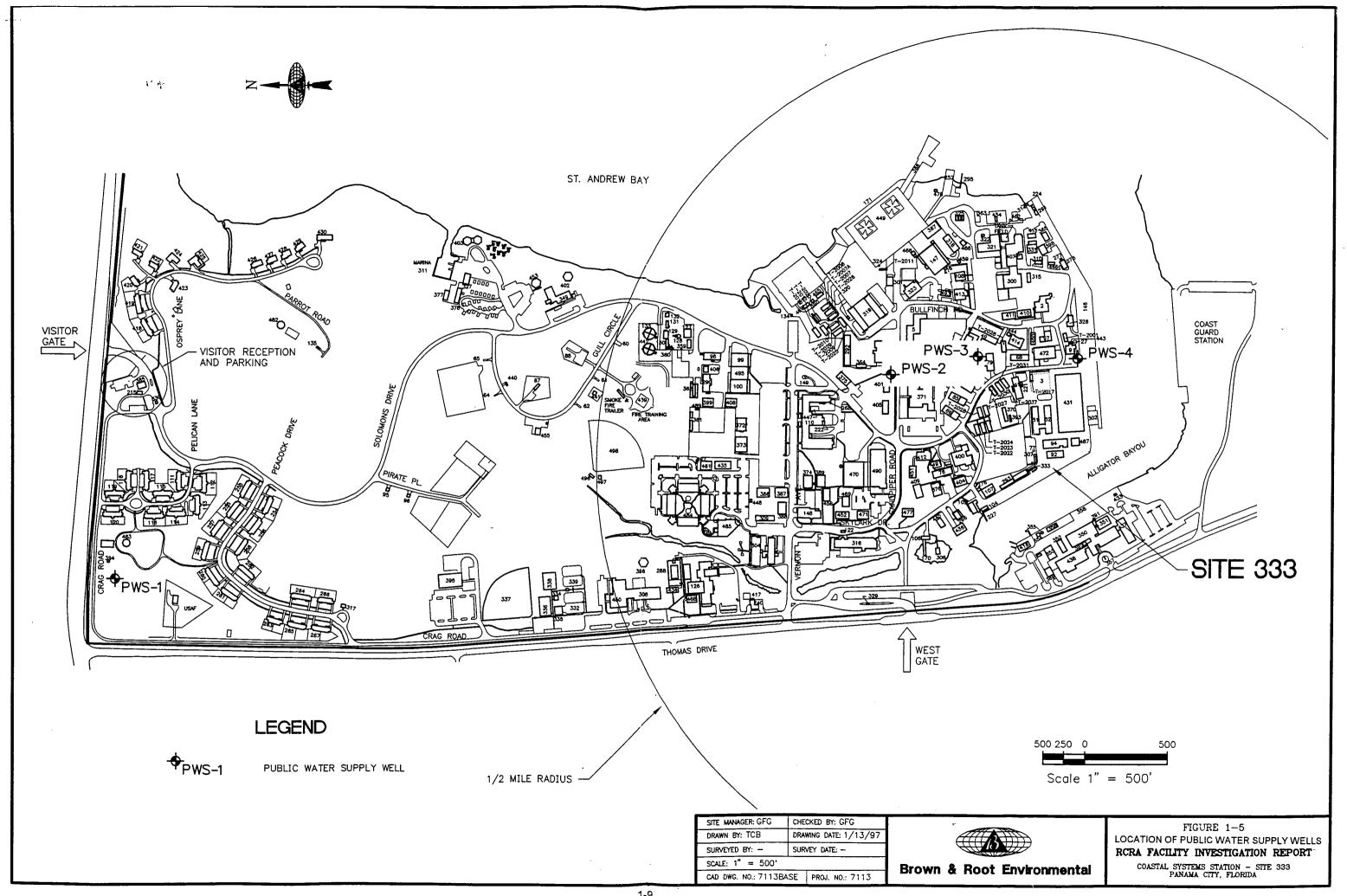
The potable water supply information presented in this report was obtained from the Resource Conservation and Recovery Act (RCRA) Facility Investigation completed for CSS (ABB Environmental Services Inc., 1995). According to this report, potable water for most of Panama City and Panama City Beach, including CSS, is supplied by surface water. Panama City Beach also uses groundwater from the Floridan aquifer system, as do private and domestic water systems throughout Bay County.

The CSS is provided potable water from the Bay County Water System, operated by the Bay County Public Utilities Department. The system draws surface water from Deer Point Lake, located 7 miles northeast of CSS. The use of county water in urban areas such as Panama City, has been reported at 83 to 95 percent.

Panama City Beach operates a public water system which uses a combination of groundwater withdrawal and surface water. The groundwater is obtained from 13 wells located in western Bay County and surface water is purchased from the county water system.

The RCRA Facility Investigation report indicates records from the Northwest Florida Water Management District list 42 permitted wells screened in the surficial aquifer system in the vicinity of CSS. These 42 wells are classified as domestic or other public supply. The permitted wells are 2-inch and 4-inch-diameter wells with capabilities generally less than 20 gallons per minute.

Four public water supply wells are located at CSS. The location of the wells are provided on Figure 1-5. These wells have 12-inch diameter casings and are completed at depths of 350 to 400 feet below land surface (bls). Of the four wells, only PWS-1, located near the housing area at Building 394 adjacent to highway 98, is currently in use. It is used to provide water for air conditioning and heat pumps only and draws water from the Floridan aquifer system at approximately 400 feet bls. The remaining wells are inactive.



No private potable wells or public potable supply wells were identified in the RCRA Facility Investigation report as being within a 1/4-mile and 1/2-mile radius of the site, respectively.

#### 1.3 SITE HISTORY AND OPERATIONS

#### 1.3.1 Site History

CSS is one of seven major research, test, and evaluation laboratories of the Space and Naval Warfare Systems Command. The site was first established in 1942 as a harbor for World War II convoy ships and as a liaison with a nearby shipyard. It later became an amphibious landing craft operations school. Research and development began in 1945 when the facility was renamed the U.S. Navy Research Countermeasures Station. In 1952 a research and development program for the use of helicopters for mine countermeasures operations was implemented at the Base. The facility was redesignated as the Naval Coastal Systems Center in 1978 and again as Coastal Systems Station in January 1992 (ABB Environmental Services, Inc. 1995).

Site 333 is utilized as a spill containment boom wash area. The site consists of a containment area, an underground waste oil tank (holding tank) and oil water/separator. The oil/water separator was primarily used to collect rinse water from diesel booms used to contain spills. Bowser-tank trucks used to haul waste oil were previously stored in the boom containment wash area. Some used oil from these trucks may have been processed through the oil/water separator system. The oil/water separator also receives water from a floor drain from Structure 307. Structure 307 is used as a hazardous drum storage area.

In November 1995, the oil/water collection system at Site 333 was upgraded. The system upgrade including the removal and replacement of a 550 gallon waste oil tank, removal of a below surface grade oil/water separator, and replacement of underground piping associated with the oil/water collection system. Prior to November 1995, a second oil/water separator with a 5 gpm capacity existed at Site 333. The separator was located south and adjacent to the above ground oil/water separator. This separator received water from the floor drain in Building 307 and from floor drains in the boom containment area. Any oils collected in the separator gravity drained into the 550-gallon fiberglass underground waste oil tank. Review of construction plans show the oil/water separator collection system which was upgraded was installed in the late 1970's (NAVFAC DRAWING No. 504-6560).

#### 1.3.2 Structural Integrity of Tanks and Lines

No structural integrity testing on the oil/water separator tank, waste oil tank, or integral piping were performed prior to removal.

#### 1.3.3 Previous Investigations

A Tank Closure Assessment was performed by Southern Waste Services (SWS) in November 1995. During removal of the oil/water separator, waste oil tank, and integral piping, SWS collected soil samples from the vadose zone soils for hydrocarbon vapor screening using an organic vapor analyzer (OVA). Four soil sampling points were established at the corners of the excavation for the waste oil tank. Three sampling points were located along the piping trench and within the oil/water separator area. Results of the soil vapor screening identified "excessively contaminated" as defined by Chapter 62-770.200 FAC., near the southeast corner of the excavation pit for the waste oil tank. Hydrocarbon vapor concentrations in the soil samples from the excavation near the holding tank ranged from less then 10 ppm to 950 ppm. Further investigation identified a greenish dark gray oily sheen within the smear zone of the water table. Soil vapor readings of soils collected along the pipe line trench and oil/water separator identified hydrocarbon vapors at concentrations of less than 10 ppm.

Two additional exploratory borings were conducted by SWS in the area between the drum containment fence and the excavation for the waste oil tank. These borings were advanced to the water table. "Excessively contaminated" soils were identified at each boring location at a depth of 4.5 feet below land surface (bls). Diesel or old hydrocarbon fuel odors were detected. A greenish gray oily soil was identified at the capillary fringe zone in one of the borings at 4.5 feet bls.

A composite soil sample for laboratory analysis was collected during the waste oil tank excavation. The sample was collected at a depth of 4 feet bls and analyzed for Volatile Organics (EPA Method 8240), Semivolatile Organics (EPA Method 8270), Total Recoverable Petroluem Hydrocarbons (TRPH) (EPA Method 9073) and RCRA Metals. In addition, soil samples were collected from the excavation wall adjacent to the oil/water separator and from the area between the drum containment and holding tank area (samples were collected at approximately 4.5 feet bls). These samples were subject to the laboratory analysis mentioned above. A composite sample was also collected from the excavated soil for lab analyses, including TCLP Volatiles and TCLP metals.

Laboratory results of the soil analyses reported Volatile Organics and TCLP Volatiles and TCLP metals below laboratory detection limits. Concentrations of RCRA metals, TRPH, 1-methynaphthalene, 2-methy naphthalene, and naphthalene were detected in the soils. Lead, cadmium, silver, and mercury were detected at 30.4 mk/kg, 1.0 mg/kg, and 0.104 mg/kg, respectively. The highest concentrations of barium and chromium were reported at 2.3 mg/kg, respectively. The highest concentrations of TRPH, 1-methy naphthalene, 2- methylnapthalene, and naphthalene were reported in soils collected near the holding tank at concentrations of 960 mg/kg, 4,800 mg/kg, 7500 mg/kg and 1,300 mg/kg, respectively.

A temporary groundwater monitoring well screened into the top of the water table was installed in the area adjacent to the holding tank. A groundwater sample was collected from the well on December 15, 1995 and analyzed for Volatile Aromatic Hydrocarbons (EPA Method 602), and Polynuclear Aromatic Hydrocarbons (EPA Method 610). Results of the sampling event report benzene, ethyl benzene, and xylenes at concentrations of 1.5 ug/L, 1.0 ug/L, and 5.2 ug/L, respectively. 1-Methynaphthalene was detected in the groundwater at 49 ug/L.

An Initial Remedial Action (IRA) Notification Form reported approximately 2 cubic yards of "excessively contaminated" soil was removed from the waste oil tank excavation during the tank closure assessment activities. The IRA Notification Form is included in Appendix A and the Tank Closure Assessment Report is provided in Appendix C.

#### 2.0 SUBSURFACE INVESTIGATION METHODS

#### 2.1 QUALITY ASSURANCE

The site investigation was conducted in accordance with the Standard Operating Procedures prescribed by the FDEP Quality Assurance Section Document DER-001/92, and adopted by the B&R Environmental Comprehensive Quality Assurance Plan Number 870055G.

#### 2.2 SOIL BORING PROCEDURES

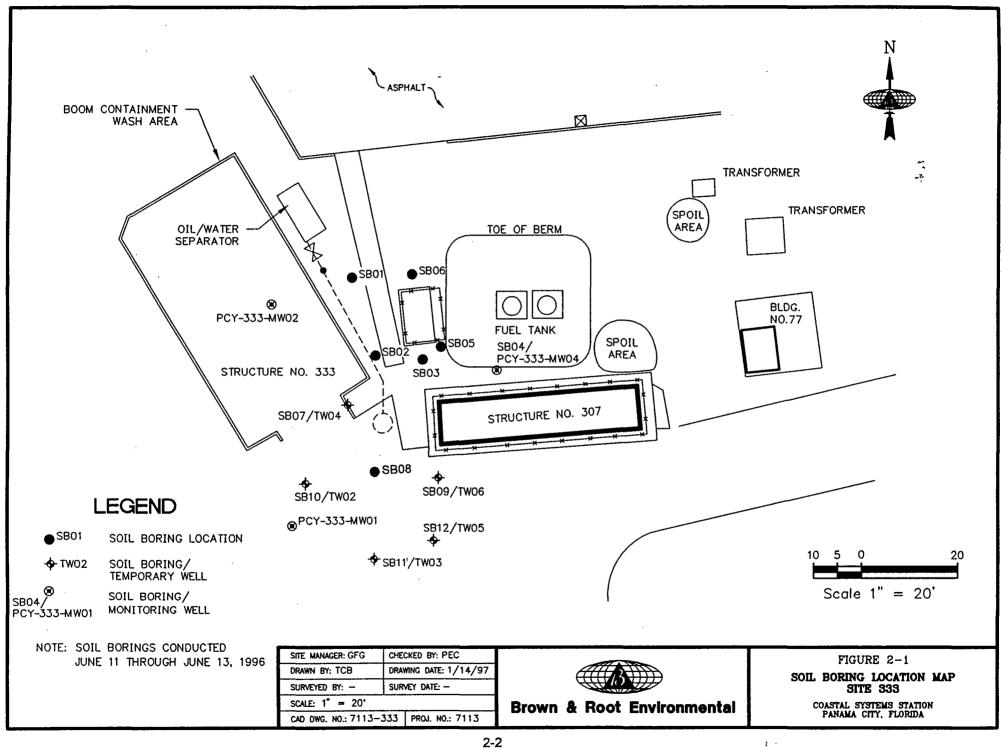
#### 2.2.1 Hand-Auguered Soil Borings

A soil hydrocarbon vapor assessment was conducted at the site by B&R Environmental on June 10 through 12, 1996. Twelve soil borings (SB01 through SB12) were excavated in the immediate area surrounding the oil/water separator and waste oil tank. Soil samples were collected from each boring for the purpose of organic vapor screening and for lithologic description. Results of the soil vapor screening would be collaborated with the soil vapor data collected during removal of the oil/water separator to determine the horizontal and vertical extent of petroleum contamination in the vadose zone. Soil borings were advanced using a 3.5 inch inside diameter (ID) stainless steel bucket auger. Soil samples were collected at two foot intervals until the water table was encountered. Wet soils were present at depths ranging from approximately 4 to 5 feet bls. Soil boring locations and soil sample collection depths are summarized on Figure 2-1 and Table 2-1, respectively. Soil boring logs are provided in Appendix D.

Prior to the advancement of the hand auger at each boring location, the hand auger was decontaminated according to B&R Environmental Comprehensive Quality Assurance Plan.

#### 2.2.2 Drilling and Soil Sampling Methods

From June 11 through June 13, 1996, borings for monitoring wells PCY-333-MW01, PCY-333-MW02, PCY-333-MW03 and PCY-333-MW04 were drilled by Groundwater Protection, Inc. under the supervision of a B&R Environmental geologist. The borings were advanced to facilitate the installation of groundwater monitoring wells. Soil samples collected during the borehole advancements were used to characterize the site lithology and provide additional assessment data on soil hydrocarbon vapor concentrations in the area. The location of the borings are shown



## TABLE 2-1 SOIL VAPOR MEASUREMENTS Coastal Systems Station Site 333

# Panama City, Florida FDEP FACILITY No. 038518667

			Headspace Readings (ppm)			
Soil Boring No.	Date of Measurement	Sample Interval (feet bis)	Total Organic Reading	Carbon Filtered Reading	Net Reading	
SB01	06-10-96	2	5	3	2	
		4	ND	ND	ND ·	
SB02	06-10-96	2	· ND	ND	ND (	
		4	10	8	2	
SB03	06-11-96	2	10	8	2	
		4	29	10	19	
SB04	06-11-96	2	2	ND	2	
(PCY-333-MW04)		4	2	ND	2	
SB05	06-11-96	2	ND	ND	ND	
		4	9	9	ND	
SB06	06-11-96	2	ND	ND	ND	
		3	.4	4	ND	
SB07	06-11-96	2	ND	-	ND	
(TW04)		4	ND	-	ND	
SB08	06-11-96	2	ND	ND	ND	
		4	ND	ND	ND	
SB09	06-11-96	2	3	2	1	
(TW06)		4	5	2	3	
SB10	6-11-96	2	ND	ND	ND	
(TW02)	1	4	ND	ND	ND	
SB11	6-11-96	2	1	ND	1	
(TW03)		4	1	ND	1	
SB12	6-12-96	2	ND	ND	ND	
(TW05)		4	5	5	ND	
PCY-333-MW01	6-12-96	2	ND	ND	ND	
		4	ND	ND	ND	
PCY-333-MW02	6-12-96	2	ND	ND	ND	
		4	8	2	6	
PCY-333-MW03	6-13-96	2	ND	ND	ND	
		4	ND	MD	ND	

#### Notes:

- = not analyzed

bls = below land surface

ppm = part per million equivalent methane

Wet soils encountered at approximately 4.5 feet bls.

on Figure 2-1, and soil boring logs are included in Appendix D.

Buried utilities were investigated at each boring location by advancing the soil boring with a post hole digger from 0 to 4 feet bls. The borings were continued with a truck mounted drill rig, using 4 1/4-inch (ID) hollow stem augers. Soil samples were collected using a split spoon sampler and standard penetrations tests were conducted in accordance with the American Society for Testing and Materials (ASTM) D-1586 recommended procedures.

Prior to the collection of the soil samples and well installations, the auger flights, drill rods, and split spoons were decontaminated according to B&R Environmental Comprehensive Quality Assurance Plan.

Soil samples were visually inspected for evidence of oil staining. Headspace analysis was conducted on each soil sample collected above the water table during the soil vapor assessment. Grab samples were collected at two foot intervals from approximately 0 to 4 feet bls. Soil vapor analysis was performed in accordance with the headspace method presented in detail in Appendix E. Hydrocarbon vapor concentrations from soil vapor analysis are summarized in Table 2-1.

Soil cuttings generated during the well installations were placed in a 55-gallon steel drum. A composite soil sample was collected from the drum and analyzed for TCLP (SW-846 1311) organics and metals, reactivity, and corrosivity. The soil will be transported for disposal by a licensed Florida waste hauler.

#### 2.3 WELL CONSTRUCTION

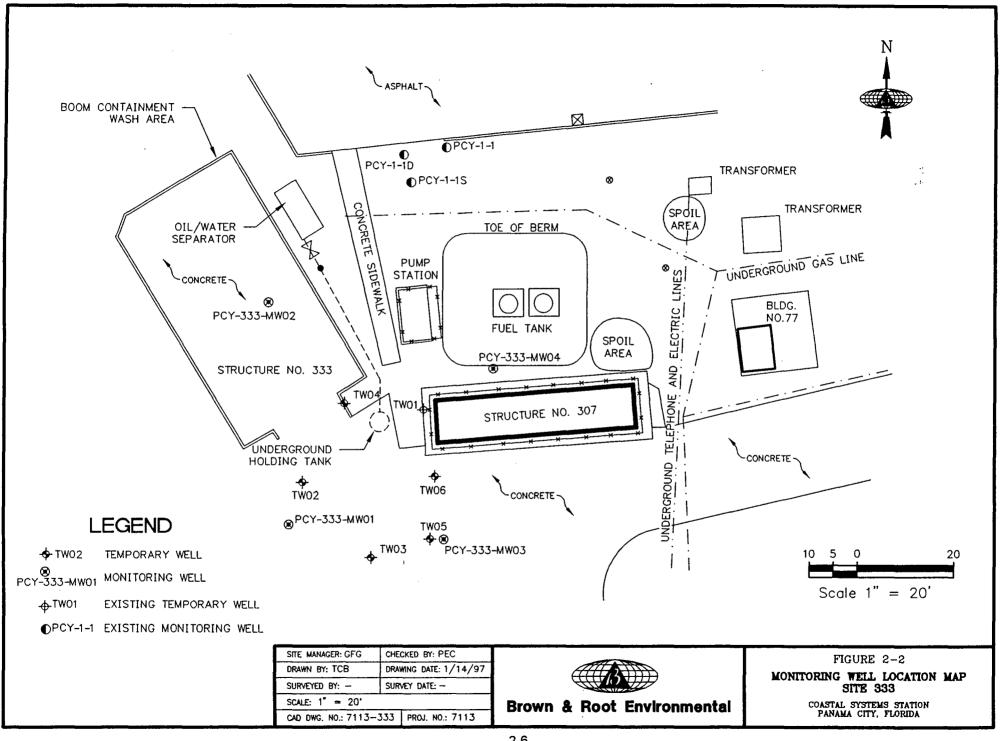
During the soil vapor assessment, temporary groundwater monitoring points TW02, TW03, TW04, TW05 and TW06 were installed in the hand augured soil borings for SB10, SB11, SB07, SB12, and SB04, respectively. These boring locations were selected for monitoring point installations based on the detection of fuel like odors in soil samples collected at the water table during the soil vapor survey. A groundwater sample was obtained from each of the temporary monitoring points using a disposable Teflon bailer. The sample was screened on-site for benzene, toluene, ethyl benzene, and xylenes, utilizing an HNU Systems Model 311D GC portable gas chromatograph (GC). Results of the GC screening were used to evaluate the optimum placement for a permanent monitoring well(s). Each temporary groundwater monitoring point was constructed of

2-inch ID threaded schedule 40 PVC solid riser and 0.010-inch slots with silt trap and well bottom cap. Each monitor point was completed with 2 feet of well screen. The monitoring point was inserted into the boring annulus and hand-pushed into the top of the water table. Upon collection of the groundwater sample, the monitoring point was pulled from the boring and a one foot plug of bentonite pellets was placed at the bottom of the boring and hydrated. The remainder of the boring was then backfilled with the soil cuttings. Each monitoring point was delivered to the site factory wrapped in plastic. Prior to installing the monitoring points, the casings and well screens were decontaminated as follows:

- Laboratory grade detergent and potable water wash
- Potable water rinse
- Isopropanol rinse
- Deionized water rinse
- Air dry

Groundwater Protection, Inc., under the supervision of a B&R Environmental geologist, installed permanent monitoring wells PCY-333-MW01 through PCY-333-MW04 on June 12 and 13, 1996. The wells were installed in conjunction with the soil boring procedures discussed above in Section 2.2.2. Each well was screened to intersect the water table and was located adjacent to the oil/water separator location. Analysis of groundwater samples collected from the well would be used to evaluate water quality in the area of the oil/water separator and underground storage holding tank. Monitoring locations are show on Figure 2-2.

The borings for monitoring well PCY-333-MW01 through PCY-333-MW03 were drilled with 4 1/4-inch (ID) hollow stem augers and a Diedrick D 120C Model drill rig. Monitoring well PCY-333-MW04 was installed using a 3.5 inch ID stainless steel bucket auger due to restricted drill rig access at location SB04. The wells were constructed of 2-inch ID, threaded, schedule 40 PVC, solid riser and 0.010-inch slots with silt trap and well bottom cap. Each well was completed at approximately 13 feet below land surface. The boring annulus was filled to approximately 1 foot above the well screen with 20/30 silica sand. A one foot layer of bentonite pellets was placed above the sand pack and hydrated. The remainder of the well annulus was grouted to within 3 inches of the top of well casing. The well was secured with a locking, water-tight cap within a 8-inch diameter steel manhole. The manhole was set within a 24-inch square concrete apron finished slightly above grade. The top of the well is secured with a locking, water-tight cap. Well completion logs are provided in Appendix F.



The wells were developed using a centrifugal pump. During well development, field measurements of pH, temperature, and specific conductance were monitored from the purge water. Each well was developed up to a maximum of one hour or until the field measurements became stable and the purge water clear. Water quality stabilization was determined using the following criteria: temperature +/-05°C, pH +/-0.1 unit, and specific conductance +/-10 umhos/cm. The wells were developed under the supervision of a B & R Environmental geologist. All development water will be removed for proper disposal by a Florida licensed waste hauler.

#### 2.4 LITHOLOGIC SAMPLING

Representative soil samples were collected to assess the shallow subsurface geologic conditions at the site. Samples used for lithologic description were collected from a stainless steel hand auger or split spoon sampler during the soil boring and monitoring well installations. Soil boring logs are included as Appendix F.

#### 2.5 SOIL VAPOR ANALYSIS

Headspace analysis was conducted on each soil sample using an Organic Vapor Analyzer-Flame Ionization Detector (OVA-FID). The soil vapor analysis was performed according to the headspace method prescribed in Rule 62-770.200 (2) FAC. Screened soil samples with corrected headspace levels in excess of 50 ppm are defined as "excessively contaminated" soil at diesel and used oil contaminated sites. The headspace methodology for determining soil organic vapor concentrations is described in detail in Appendix E.

#### 2.6 SOIL SAMPLE ANALYSIS

Analysis of soil samples collected during the Tank Closure Assessment were used to determine the presence or absence of petroleum hydrocarbons or waste oil constituents in the area of the oil/water separator and holding tank. No additional subsurface soil samples were collected for laboratory analysis during the contamination assessment investigation. Soil sample collection and parameters analyzed during the Tank Closure Assessment are discussed in Section 1.3.3. Soil quality results from the Tank Closure are included in Appendix C.

#### 2.7 HYDROLOGIC INVESTIGATION

#### 2.7.1 Water Level Measurements

The depth to water was measured in existing site wells PCY-1-1S and TW01, and in monitoring wells PCY-333-MW01 through PCY-333-MW04, on July 11, 1996 and November 25, 1996. Monitoring well PCY-1-1S was installed as part of the Resource Conservation and Recovery Act Facility Investigation for SWMU-1 (ABB Environmental Services Inc., 1995). This well is a 2-inch ID PVC well screened to intersect the water. The well terminates at approximately 13 feet bls. Existing temporary well TW01 was installed during the Tank Closure Assessment completed for Site 333. TW01 is constructed of 2-inch ID PVC and is screened from 3 to 8 feet bls.

Water level measurements were collected from the top of well casings using an electronic water level indicator. Each top of well casing was surveyed to the National Geodetic Vertical Datum (NGVD) 1929 by a Florida licensed Professional Land Surveyor. Subtracting the depth to water level measurements from the well top casing elevations provided the water table elevation relative to the NGVD. The water level field forms and field elevation survey data are included in Appendix H. The water table gradient across the site was evaluated from water level measurements collected on July 11, 1996. The groundwater gradient was calculated by determining the perpendicular distance between groundwater contours developed from groundwater elevation data. Gradient calculations are included in Appendix G. The hydraulic conductivity value obtained from the RCRA Field Investigation study for SWMU-1 was used in calculating the site's groundwater velocity. The effects of tidal influence on the site was evaluated from tidal survey data collected for the RCRA Field Investigation for SWMU-1.

#### 2.7.2 Aquifer Characteristics

The hydraulic conductivity value obtained from the RCRA Field Investigation study for SWMU-1 was used to evaluate the hydraulic conductivity and groundwater velocity for the site.

#### 2.7.3 Groundwater Flow Velocity and Transmissivity

The groundwater flow gradient was evaluated using the following equation:

$$i = (h_1 - h_2)/d$$

where:

i = the hydraulic gradient

h<sub>1</sub> = the water elevation at point 1

h<sub>2</sub> = the water elevation at point 2

d = the distance between point 1 and point 2

Potential movement of groundwater at the site may be described in terms of transportation by natural flow system in the saturated zone while assuming groundwater flow follows Darcy's Law. Darcy's Law may be expressed as:

$$V = (K^*i)$$

$$n_e$$

where:

V = average velocity K = hydraulic conductivity i = hydraulic gradient n<sub>e</sub> = effective porosity

Site specific transmissivity is calculated using the following equation:

$$T = K*b_e$$

where:

T = transmissivity

K = hydraulic conductivity

be = effective aquifer thickness

The groundwater flow velocity and aquifer transmissivity calculations are included in Appendix G.

#### 2.7.4 Tidal Influence Survey

A tidal survey was conducted during the RCRA Facility Investigation to determine if the hydraulic gradient at locations close to Alligator Bayou is influenced by tidal fluctuations. Continuous water level measurements were obtained from several selected monitoring wells for a period of 24 hours. Monitoring wells PCY-14-5 and PCY-1-3 were selected at SWMU 1 and Area of Concern (AOC) 2, to evaluate the effects of tidal influence by Alligator Bayou. Monitoring well PCY-14-5 is located 40 feet from the seawall at Alligator Bayou and was paired with PCY-1-3, located 200 feet from the Bayou.

#### 2.8 WATER SAMPLING

#### 2.8.1 Free Product Sampling

Prior to groundwater sampling, B&R Environmental personnel checked each well for free product using a pre-cleaned Teflon® bailer. The Teflon® bailer was used to extract a water sample from the top of the well's water column to visually inspect for free product. Oil films (films less then 0.01 foot thick) were observed in wells PCY-333-MW04 and TW01 on July 11, 1996. A free product thickness of 0.15 feet was detected in the well PCY-333-MW04 on November 25, 1996.

#### 2.8.2 Groundwater Sampling

Groundwater sampling was performed to determine the presence or absence of dissolved petroleum hydrocarbons or used oil constituents in shallow groundwater in the vicinity of the former oil/water separator. Groundwater samples were collected by B&R Environmental personnel from wells PCY-333-MW01, PCY-333-MW02, PCY-333-MW03, and PCY-1-1S on July 11, 1996. Groundwater samples were analyzed by ICP Series for lead, arsenic, cadmium, and chromium (all metal analysis were unfiltered), EPA Method 504.1 for Extractable Volatile Organic (1-2dibromoethane EDB), EPA Method 601 for Purgeable Halocarbons, EPA Method 602 for Purgeable Aromatics (benzene, toluene, ethyl benzene, and methyl-tert butyl either). EPA Method 610 for PAHs, SW-846 Method 8260 for GC/MS Volatile Organics, SW-8468 Method 8270A for GC/MS Semivolatile Organics and EPA Method 418.1 for Total Petroleum Hydrocarbons. The lead, arsenic, and chromium samples were collected using new silicon tubing and a peristaltic pump. The remainder of the groundwater parameters analyzed during the sampling event were collected using pre-cleaned Teflon® bailers. Approximately five well volumes of groundwater were removed from each well using a pre-cleaned Teflon® bailer. Temperature, pH, specific conductance measurements and well purge volumes were recorded at the time of sample collection and are provided in Appendix H. Groundwater samples were placed on ice and shipped via overnight courier to Quality Analytical Laboratories, Inc., Montgomery, Alabama.

Groundwater samples were not collected from wells PCY-333-MW04 and TW01 during the July 11 sampling event due to the presence of an oil film in the water at the well locations. The oil films were identified during visual inspection of the water clarity using a Teflon® bailer. On November 25, 1996, groundwater samples were collected from wells PCY-333-MW04 and TW01 for the test parameters, as referenced above. During the November 25 sampling event, no product film was observed on the water in TW01. A free product thickness of 0.15 feet was measured in PCY-333-

MW04. Prior to sample collection, the product in the well was bailed and removed from the well. An unfiltered lead sample was also collected from PCY-333-MW02 during the November sampling event to provide additional data to evaluate lead concentrations in the area near PCY-333-MW02.

Groundwater samples were collected in accordance with the FDEP Quality Assurance Document DER-001/92. During the sampling events, quality control samples (i.e. equipment blanks, trip blanks, duplicate) were prepared and submitted to the laboratory as required by the approved QA procedures. Sampling activities were documented in a site specific field logbook, and samples were transmitted under chain-of-custody protocols.

#### 3.0 RESULTS OF INVESTIGATION

#### 3.1 SITE HYDROGEOLOGY

#### 3.1.1 Lithology

Lithologic samples collected by B&R Environmental indicate the site is predominantly underlain by dark-gray, to light-tan, fine to medium grained sand and silty sand to a depth of approximately 13 feet bls. Due to the homogeneity of the subsurface, no lithologic cross-section was constructed. Soil boring logs are included as Appendix E.

The RCRA Facility Investigation (ABB Environmental Services, Inc.,1995) indicates sand is the primary soil type encountered to a depth of 27 feet bls at well PCY-1-1D, located approximately 40 feet northeast of the oil/water separator. A dark green sandy clay and clayey sand is present at this locality between 27 and 35 feet bls. The sand, sandy clay, and clayey sand deposits in the surficial aquifer are underlain by olive-green, fossiliferous, clayey limestone of the Intracoastal Formation at a depth of 48 feet bls.

#### 3.1.2 Aquifer Characteristics and Classification

Based on water level data collected from site monitoring wells on July 11 and November 25, 1996, the depth to the shallow aquifer at the site is approximately 4 to 5 feet bls. The groundwater level measurements are presented in Table 3-1. The water level measurement field forms are provided in Appendix H. The aquifer is classified as a G-II aquifer based on dissolved solids content typically associated with the surficial aquifer in the area of CSS.

Rising head slug tests of the surficial aquifer at CSS for the area of SWMU-1were conducted to estimate the hydraulic conductivity during the RCRA Facility Investigation. The geometric mean hydraulic conductivity for the surficial aquifer for SWMU-1 was estimated at  $4.98 \times 10^{-3}$  feet/minute (7.1 ft/day) (ABB Environmental Services Inc., 1995). Due to the proximity of Site 333 to SWMU-1, the hydraulic conductivity value determined for SWMU-1 was used for Site 333.

Using the groundwater flow gradient equation presented in Section 2.7.3, a hydraulic gradient of 0.05 feet/foot to the south-southwest was calculated from the data collected on July 11, 1996; these data are represented in Figure 3-1. The November 25, 1996, water level data indicates a

#### **TABLE 3-1 DEPTH TO GROUNDWATER MEASUREMENTS Site 333**

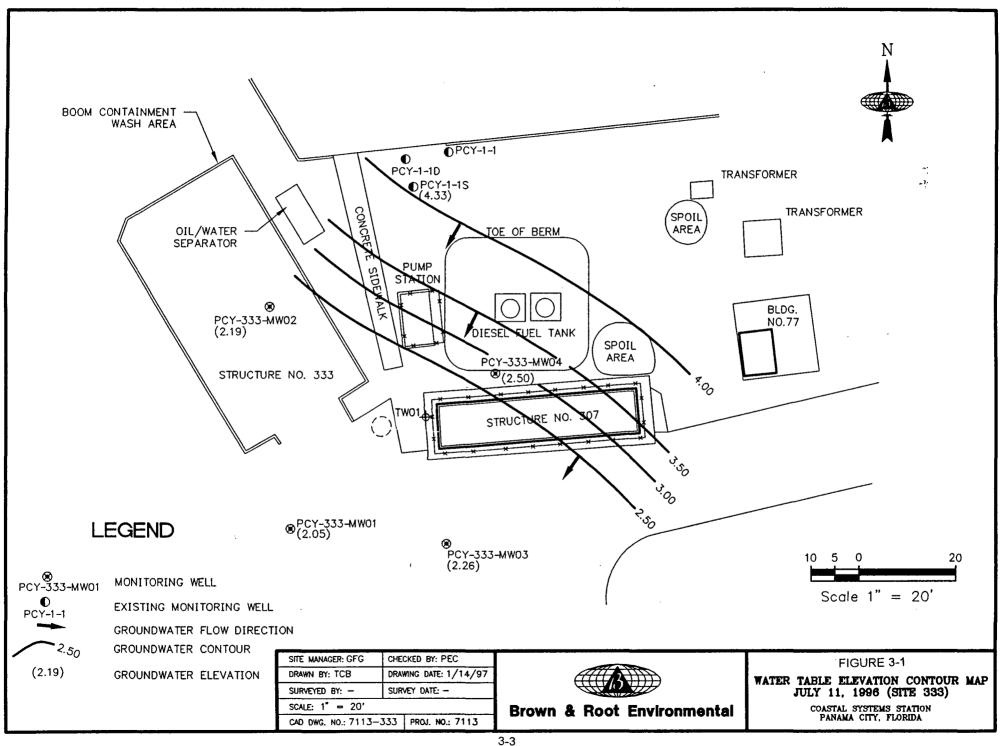
## Coastal Systems Station, Panama City, Florida FDEP Facility No. 038518667

Monitoring Well ID	Date	Top of Well Casing Elevation (feet NGVD)	Free Product Thickness (feet)	Depth to Water (feet bls)	Water Table Elevation (feet NGVD)	Well Screen Interval (feet bis)
PCY-333-MW01	07/11/96	6.65	0.00	4.60	2.05	3 to 13
	11/25/96	6.65	0.00	4.77	1.88	
PCY-333-MW01	07/11/96	6.39	0.00	4.20	2.19	3 to13
	11/25/96	6.39	0.00	4.40	1.96	
PCY-333-MW-03	07/11/96	6.57	0.00	4.31	2.26	3 to 13
	11/25/96	6.57	0.00	4.58	1.99	
PCY-333-MW04	07/11/96	7.75	<.01	5.25	2.50	3 to 13
	11/25/96	7.75	0.15	5.80	1.95	
PCY-1-1S	07/11/96	9.48	0.00	5.15	4.33	2 to 10
	11/25/96	9.48	0.00	6.60	2.88	
TW01	07/11/96	8.70	<0.01	5.44	3.26	5 to 15
i 	11/25/96	8.70	0.00	6.76	1.94	

Notes:

bls = below land surface. ID = identification

NGVD = elevation relative to the National Geodetic Vertical Datum 1929.



much flattter flow gradient (.002 feet/foot) toward the south-southwest as shown on Figure 3-2.

Lithologic data and available literature indicate the effective porosity of the sediments comprising the surficial aquifer is approximately 0.30 (Heath, 1994).

Using a hydraulic conductivity of 7.1 feet/day, the hydraulic gradient of 0.05 feet/foot, an inferred effective porosity value of 0.30, and Darcy's Equation as stated in Section 2.7.3, the groundwater flow velocity across the site is calculated at 1.81 feet/day in a south-southwest direction. The transmissivity of the surficial aquifer was calculated at 156 ft<sup>2</sup>/day.

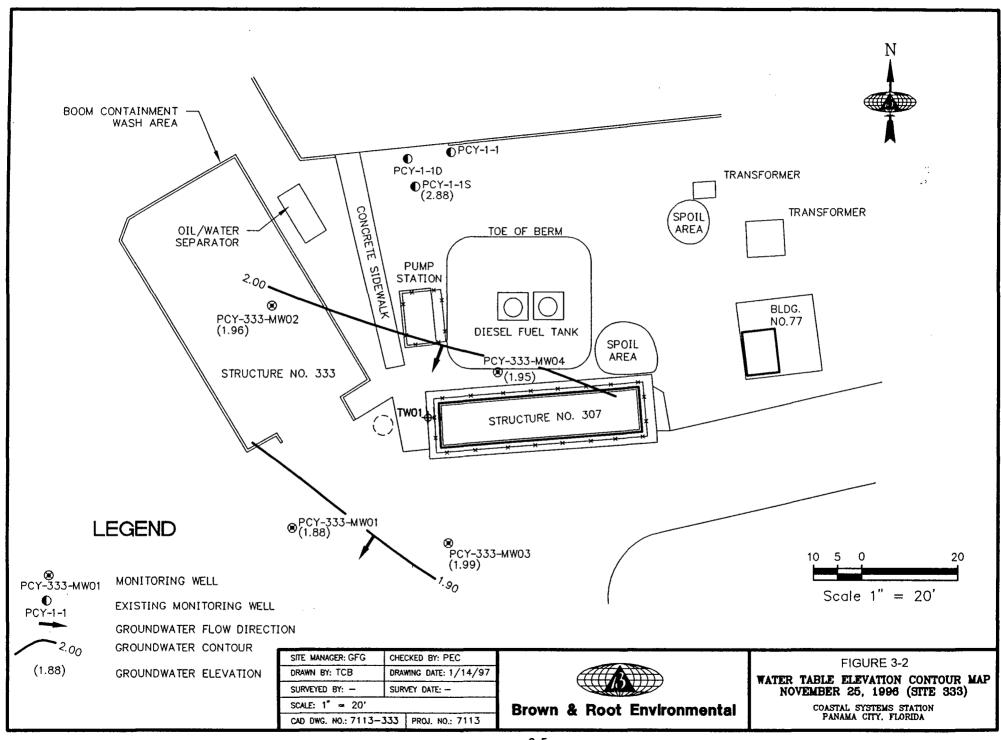
#### 3.1.3 Tidal Influence

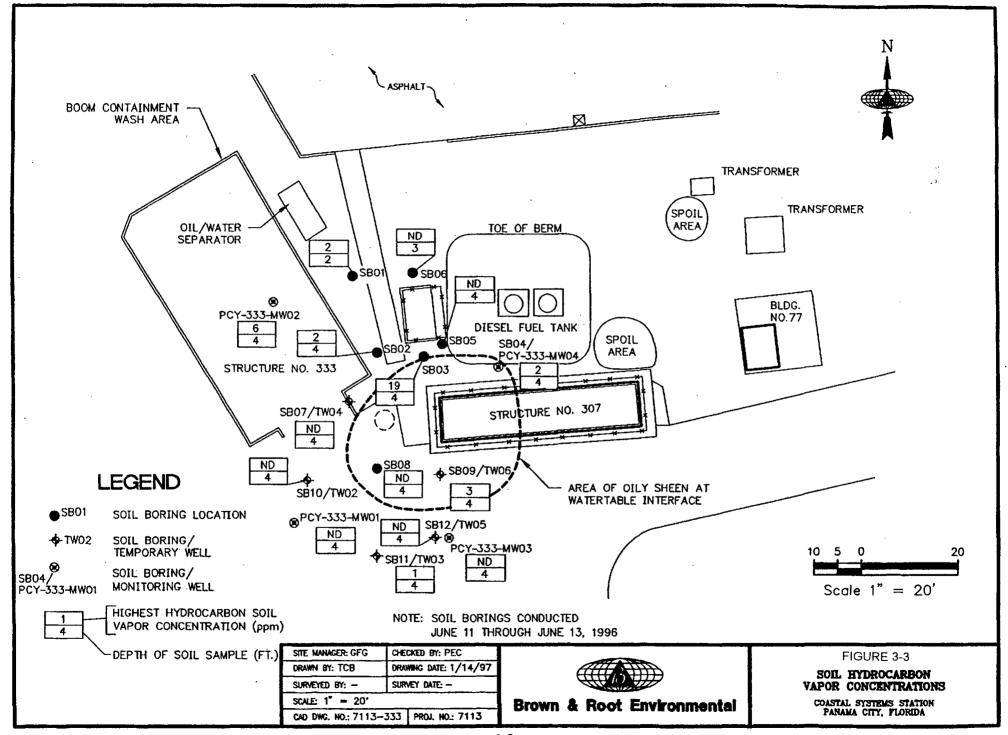
Results of the tidal survey conducted at monitoring wells PCY-14-5 and PCY-1-3 indicate groundwater flucuations due to tidal influence is limited to a distance of less than 200 feet at SWMU 1 on Alligator Bayou. The distance in which significant tidal influence (that is, influence strong enough to affect groundwater flow) effects groundwater elevation is less than 40 feet from the bayou. This may be due to the presence of the sea wall at the bayou which may interfere with natural groundwater flow (ABB Environmental Services Inc., 1995).

#### 3.2 SOIL QUALITY

The highest soil hydrocarbon vapor concentrations detected in vadose zone soils were 19 ppm and 6 ppm at boring locations SB03 and PCY-333-MW02, respectively. These samples were collected at a depth of 4 feet bls at the water table interface. No "excessively contaminated" soil as defined by Chapter 62-770.200, FAC was encountered in any of the soil samples. Soil vapor screening results are presented in Table 2-1. Soil boring locations and vapor readings are depicted on Figure 3-3.

Soil samples collected from the saturated zone at 4.5 feet to 5 feet bls at borings SB04, SB06, SB07, SB08, and SB09 exhibited an oil sheen with a fuel like odor. Soil analyses conducted on the soil at depths of 4 to 4.5 feet bls near the waste oil tank during the Tank Closure Assessment reported no volatile organic. Concentrations of 1-methynaphthalene, 2- methylnaphthalene, naphthalene, and TRPH were detected in the soil which are compounds characteristic of diesel fuel. No TCLP Volatiles or TCLP metals were detected in the composite soil sample collected from the "excessively contaminated soil". Soil laboratory analytical data sheets are included in the





Tank Closure Assessment provided in Appendix C.

### 3.3 WATER QUALITY

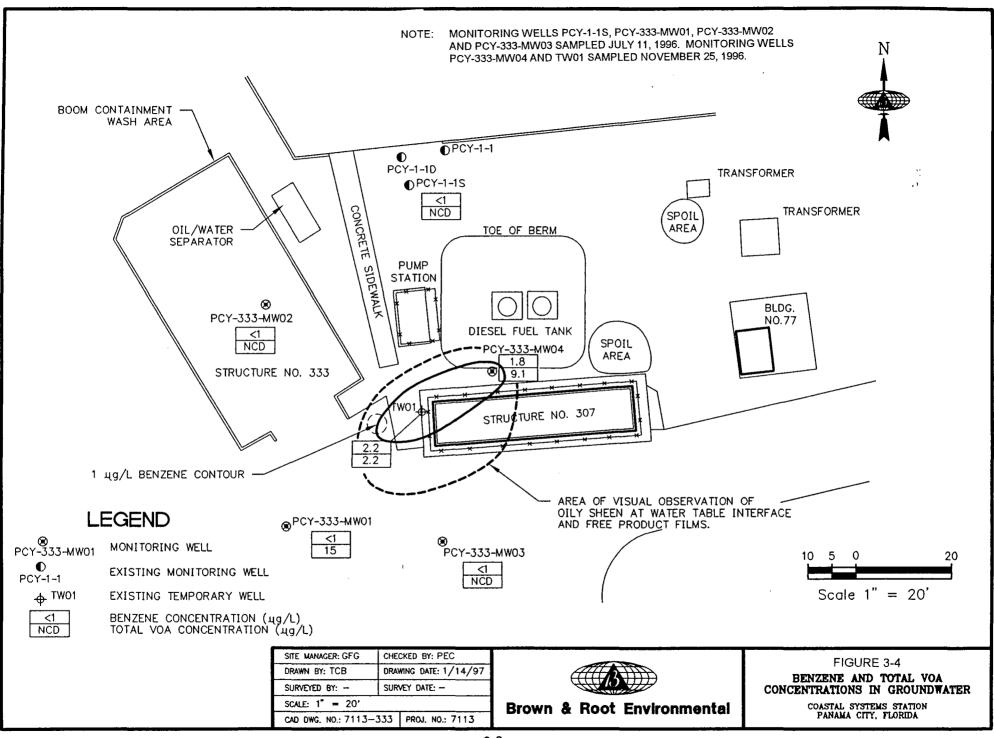
Results of field screening groundwater samples with a portable GC identified dissolved hydrocarbons at TW02 (SB10) and TW06 (SB11). The groundwater sample from TW02 reported toluene, ethylbenzene, and o-xylenes at concentrations of 3, 12, and 30 parts per billion (ppb), respectively. Groundwater in well TW06 contained benzene, toluene, ethyl- benzene, m,p-xylenes, and o-xylenes at concentrations of 14, 23, 136, 43, and 169 ppb, respectively. Dissolved hydrocarbons were not detected in the groundwater sample collected form TW03 (SB11), TW04 (SB07), and TW05 (SB12). Based on the field screening water quality results, the locations for present monitoring wells PCY-333-MW01 and PCY-333-MW03 were selected to characterize groundwater quality near TW02 and TW06. TW02 and TW06 are located south and southwest of the oil/water separator. The field GC groundwater quality reports are provided in Appendix I.

Water quality results of groundwater samples collected from site monitoring wells on July 11, 1996 and November 25, 1996 are summarized as follows:

Benzene was reported at 1.8 and 2.2 microgram per liter (ug/L) in monitoring wells PCY-333-MW04 and TW01, respectively. Benzene was reported below laboratory detection limits in all other groundwater samples. The concentrations of benzene reported in groundwater samples are slightly above the State Target Level 1 ug/L. Total VOA concentrations of 15 ug/L, 9.1 ug/L, and 2.2 ug/l were reported in wells PCY-333-MW02, PCY-333-MW04, and TW01, respectively. The concentrations of Total VOAs reported in groundwater samples are well below the State Action Level of 50 ug/L. Benzene and Total VOA groundwater concentrations are provided on Figure 3-4.

Volatile Organic Halocarbons were reported in groundwater samples collected in PCY-333-MW01 and TW01. Cis-1-2-dichloroethane was detected in PCY-333-MW01 at 1.8 ug/L. This concentrations is below the Florida Primary Drinking Water Standard of 70 ug/L. Vinyl chloride was detected in TW01 at 1.4 ug/L. The Florida Primary Drinking Water Standard is 1.0 ug/L.

1,2-Dichloroethane, 1,2-dibromoethane (EDB), and volatile organics were reported below laboratory detection in all groundwater monitoring well samples.



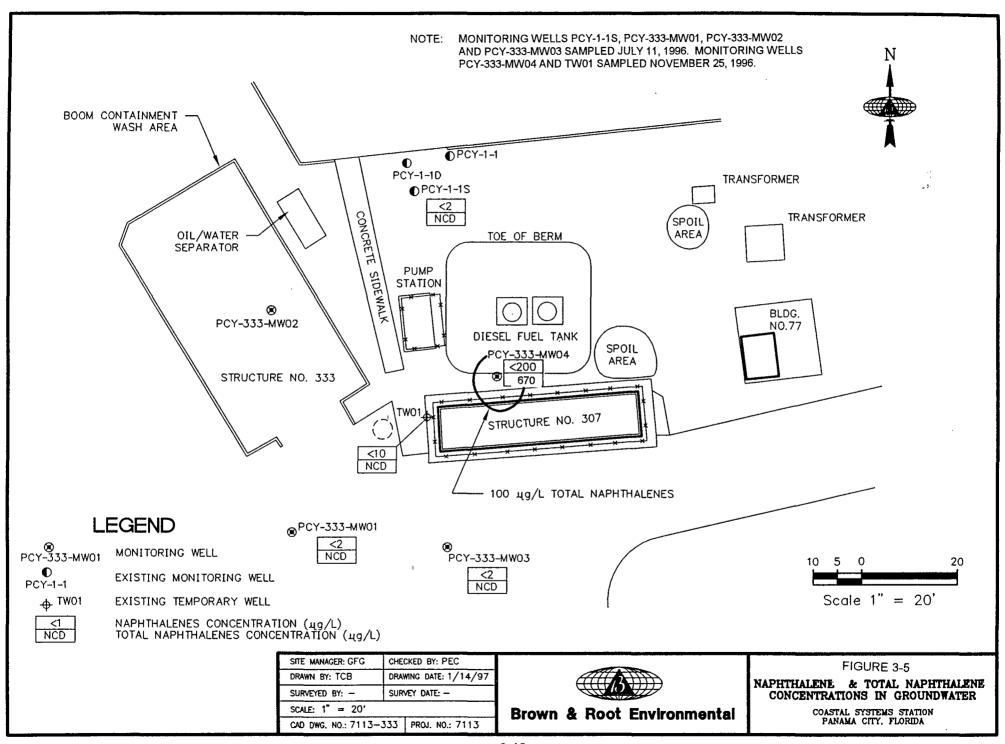
Naphthalene was reported below detection limits in all monitoring well samples. An elevated laboratory detection limit of <200 ug/l was used in the sample reporting of groundwater analysis from PCY-333-MW04. The reported total naphthalene concentration of 670 ug/L reported in this well maybe less than the actual total naphthalene concentration due to the elevated naphthalene detection limit. Prior to sampling well PCY-333-MW04, the well contained a free product thickness of 0.15 feet thickness of free product. Well PCY-333-MW04 is located adjacent to a diesel underground storage tank system (UST). Total Naphthalene was reported below detection limits in all other groundwater samples. The State Target Level for Total Naphthalene is 100 ug/l. Naphthalene and Total Naphthalene groundwater concentrations are shown on Figure 3-5.

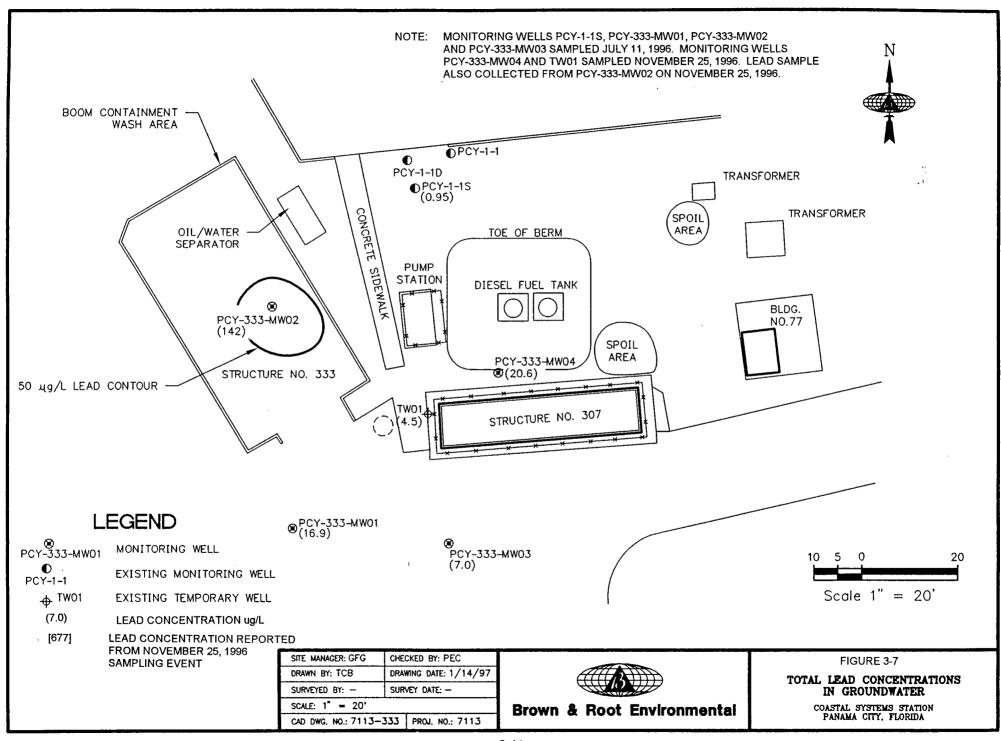
Semi volatile organics in groundwater were reported below laboratory detection limits in all monitoring well samples except PCY-333-MW04. Phenanthrene was reported in PCY-333-MW04 at 160 ug/L. Elevated laboratory detection limits were used in the reporting of groundwater quality results for PCY-333-MW04. The laboratory detection limits for the compounds analyzed in this sample generally ranged from <100 to <500 ug/L. Due to the elevated detection limits, other semi volatile organic compounds could exist in the groundwater at PCY-333-MW04 but were not detected. Phenanthrene is classified as a systemic toxicant under Florida Ground Water Guidance Concentrations. The guidance concentrations for Phenanthrene is 10 ug/L.

Arsenic and chromium were detected in several groundwater monitoring samples. The highest concentration of arsenic was detected in PCY-333-MW03 at 4.0 ug/L. The highest concentration of chromium was reported in PCY-333-MW01 at 21.4. The State Target Levels for arsenic and chromium are both 50 ug/L. Cadmium was reported below laboratory detection limits in all groundwater monitoring well samples.

Groundwater lead concentrations were reported below the State Action Level of 50 ug/L in all wells except PCY-333-MW02. Concentrations of lead were detected in PCY-333-MW03 at 142 and 677 ug/L during the July and November sampling events. PCY-333-MW02 is located within the landfill area of SWMU 1. During installation of the well, rusted metal objects were encountered at the water table which could be contributing to the elevated lead concentrations at the well location. Lead concentrations in groundwater samples collected in other monitoring wells at Site 333 ranged from 0.95 to 20.6 ug/L, as shown on Figure 3-6.

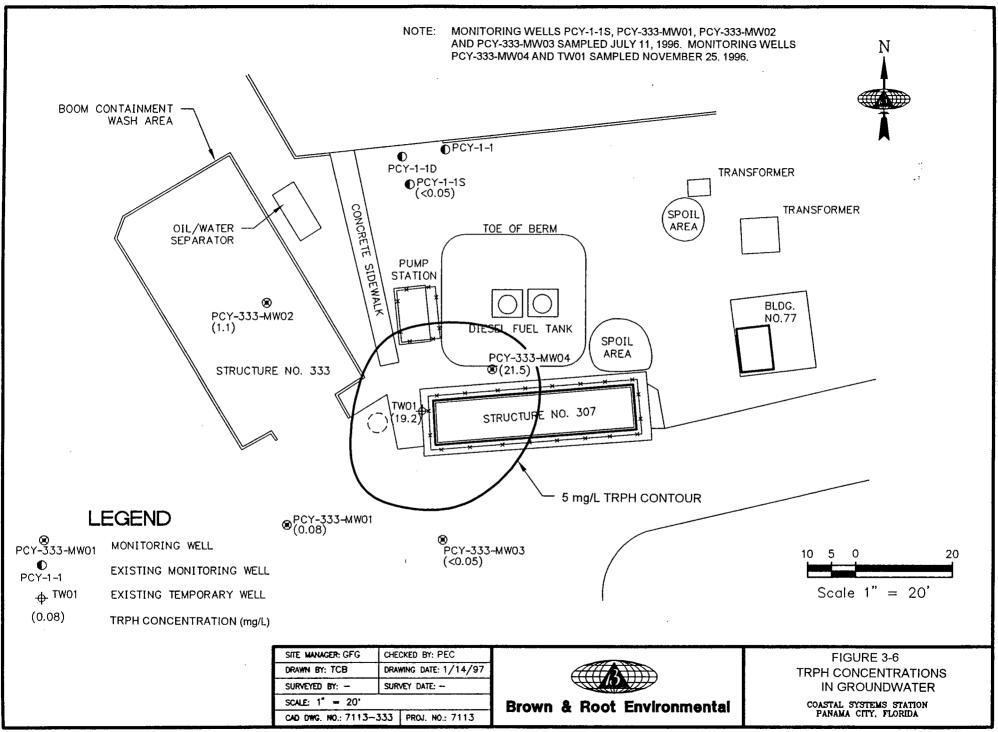
TRPH concentrations were reported in several wells. The highest concentrations were reported in groundwater samples from TW01 and PCY-333-MW04 at concentrations of 19.2 and 21.5





milligrams per liter (mg/L). These concentrations exceed the State Action level of 5 mg/L. TRPH groundwater concentrations are provided in Figure 3-7.

Toluene was reported in the Equipment Blank sample at 1.4 ug/L for samples collected on July 7, 1996. The detection of this constituents was at a level which did not effect the interpretation of the data and the water quality data is considered acceptable. The Equipment Blank sample collected on November 25, 1996 reported trace amounts of MTBE and lead. This constituents were reported at levels were which did not effect the interpretation of the data. A summary of groundwater analytical results are presented in Table 3-2. Groundwater laboratory analytical results are provided as Appendix J. Field sampling forms are included in Appendix H.



# TABLE 3-2 SUMMARY OF GROUNDWATER QUALITY: SELECTED PARAMETERS FROM THE GASOLINE AND KEROSENE ANALYTICAL GROUP

### **Site 333**

### Coastal Systems Station, Panama City, Florida FDEP ID No. 038518667

Monitoring Well ID	Date Sampled	Benzene (μg/L)	Total VOA (μg/L)	MTBE (μg/L)	DCE (μg/L)	EDB (µg/L)	NAP (μg/L)	Total NAPS (μg/L)	TRPH (mg/L)	Volatile Organics (μg/L)	Semi Volatile Organics (μg/L)	Arsenic Unfiltered Samples (μg/L)	Cadmium Unfiltered Samples (μg/L)	Chromium Unfiltered Samples (μg/L)	Lead Unfiltered Samples (μg/L)
PCY-1-1S	7/11/96	<1.0	<1.0	<1.0	<1.0	<0.02	<2	NCD	<0.05	NCD	NCD	2.2	<3.3	<2.2	0.95
PCY-333- MW01	7/11/96	<1.0	15	9.0	<1.0 *(1)	<0.02	<2	NCD	0.08	NCD	NCD	2.9	<3.3	21.4	16.9
PCY-333- MW02	7/11/96 11/25/96	<1.0 NA	<1.0 NA	<1.0 NA	<1.0 NA	<0.02 NA	<2 NA	NCD NA	1.1 NA	NCD NA	NCD NA	3.7 NA	<3.3 NA	11.6 NA	142 677
PCY-333- MW03	7/11/96	<1.0	<1.0	<1.0	<1.0		<2	NCD	<0.05	NCD	NCD	2.6	<3.3	7.5	5.8
PCY-333- MW03 (Duplicate Sample)	7/11/96	<1.0	<1.0	<1.0	<1.0	<0.02	<2	NCD	<0.05	NCD	NA	4.0	<3.3	8.9	7.0
PCY-333-	7/11/96	NA	NA	NA	NA 11.0	NA	NA	NA 070	NA 31.5	NA	NA 160 <sup>(2)</sup>	NA 14.0	NA 10.0	NA	NA
MW04	11/25/96	1.8	9.1	<1.0	<1.0	<0.02	<200	670	21.5	NCD_		<1.6	<3.3	3.0	20.6
TW01	7/11/96 11/25/95	NA 2.2	NA 2.2	NA <1.0	NA ∗(3)	NA <0.02	NA <10	NA NCD	NA 19.2	NA NCD	NA NCD	NA 1.8	NA <3.3	NA <2.2	NA 4.5
Trip Blank	7/11/96 11/25/96	<1.0 NA	1.8 NA	<1.0 NA	<1.0 NA	NA NA	NA NA	NA NA	NA NA	NCD NA	NA NA	NA NA	NA NA	NA NA	NA NA
Equipment Blank	7/11/96 11/25/96	<1.0 <1.0	1.4 <1.0	<1.0 1.3	<1.0 <1.0	<0.02 <0.02	<2 <2	NCD NCD	<0.05 <0.06	NCD NCD	NCD NCD	<0.95 <1.6	<3.3 <3.3	<2.2 <2.2	<0.75 1.4

Notes: (1)

= cis-1,2-dichloroethene detected at 1.8 ug/L

(2) = phenanthrene detected at 160 ug/L (elevated detection limit)
(3) = vinyl chloride detected at 1.4 ug/L (DEC reported at<1.0 ug/L

ID = identification ug/L = micrograms per liter NA = not analyzed

Total VOA = total volatile organic aromatics = sum of benzene, toluene, ethylbenzene, and xylenes

MTBE = methyl tert-butyl ether DCE = 1,2-dichloroethane

EDB = 1,2-dibromoethane = ethylene dibromide

NCD = no constituents detected
TRPH = total petroluem hydrocarbons

NAP = naphthalene

Total NAPS = sum of total naphthalene constituents detected in sample

### 4.0 DISCUSSION

In November 1995, approximately 2 cubic yards of "excessively contaminated" soil was removed during an oil/water system upgrade for the boom containment wash area. The "excessively contaminated" soil was from an excavation to remove the waste oil tank used as part of the oil/water system collection system. During June 1996, B&R Environmental conducted a soil vapor survey in the area of the former oil/water separator. "Excessively contaminated" soil was not encountered within the vadose zone soils during the assessment. The removal of the "excessively contaminated" soil from the vadose zone, appears to have been completed during excavation of soils in November 1995.

Samples of the oil stained soil collected during the Tank Closure Assessment reported no volatile organics in the soil. Semi volatile organic constituents with diesel fuel composition (naphthalene, 1-methylnaphthalene, and 2-methynaphthalene) were reported in the soil sample collected at 4.5 feet bls. The presence of these constituents suggests soil contamination is attributed to a diesel fuel release. TRPH was also detected in the soil analyses indicating the source of contamination is petroleum in nature. No TCLP volatiles or TCLP metals were detected in the composite soil sample collected from the "excessively contaminated" soil pile to indicate the soil is hazardous.

During the soil vapor assessment, oil stained soils were identified at several boring locations near the waste oil tank. The oil stained soils were identified in soil samples collected at the water table interface at 4.5 to 5 feet bls. Oily saturated soil and free product was also detected at monitoring well location PCY-333-MW04. Well PCY-333-MW04 is located upgradient of groundwater flow to the waste oil tank and downgradient of flow to the diesel underground storage tanks which are located approximately 25 feet northeast of the waste oil tank. It is therefore unlikely that the contamination in well PCY-333-MW04 is unrelated to Site 333.

Laboratory analysis of groundwater samples collected during the CA indicate dissolved hydrocarbon concentrations above State Action Levels for benzene, total VOAs, naphthalene, total naphthalene, TRPH, and lead. The benzene, total VOAs, naphthalene, total naphthalene, and TRPH dissolved plume concentrations appear to be limited to the southeast area of the diesel UST, and the area where excessively contaminated soils were excavated during the tank closure assessment. These plume areas also correlate with the area where visual observation of oil stained soils were identified.

The oil stained soils at the water table interface appear to be from old spills which occurred at the diesel UST. The diesel UST is located directly upgradient of the underground waste oil tank. The lateral dispersion of the oil-stained soils at the water table interface, and the low concentrations of dissolved hydrocarbons detected in groundwater samples collected from site monitoring wells (other than source well PCY-333-MW04), suggests the product is heavily weathered.

Vinyl chloride and cis-1,2,-dichloroethane were detected at concentrations slightly above laboratory detection limits. The source for these constituents is unknown and is likely unrelated to Site 333

Monitoring well PCY-333- MW02 reported elevated levels of lead above State Action Levels. The elevated lead at this location may have been caused by the disposal of scrap metal in SWMU-1. Rusted objects were encountered during the installation of the well which may have elevated the lead concentration in that well.

Based on petroleum hydrocarbon concentrations detected in monitoring wells downgradient of the site, Alligator Bayou does not appear to be threatened from the levels of hydrocarbons detected at the site.

The predominant soil type of the surficial aquifer is composed of sand. A sandy clay and clayey sand deposit was identified at a depth of 27 to 35 feet bls at monitoring well location PCY-1-1D. Well PCY-1-1D is located approximately 50 feet north of the waste oil tank. These fine grained deposits would have lower permeability then the overlying sand and should help attenuate the downward migration of petroleum hydrocarbons.

Depth to water to the surficial aquifer (water table) has been documented at approximately 4.5 to 5 feet bls. Subsurface utilities identified at the site which could potentially intersect the water table include a water main and sanitary sewer line. These utilities are typically completed at depths of 4 feet bls. These utilities are located up gradient of the dissolved contaminant plume identified at the waste oil tank and diesel tank area and should not provide a pathway for the transport of dissolved constituents in the subsurface. The direction of groundwater flow for the surficial aquifer is towards the southwest toward Alligator Bayou. The groundwater flow velocity was calculated at 1.18 feet/year.

The effects of tidal influence on the groundwater flow direction at the site is negligible. A sea wall constructed at Alligator Bayou restricts the natural flow of groundwater in the area and limits the effects of tidal influence. Alligator Bayou acts as the natural discharge point for the water table aquifer downgradient of the site. However, the seawall does create a vertical component of groundwater flow as groundwater which normally discharges to the bayou tries to flow under and/or along the seawall.

The well fields and surface water intakes which supply drinking water to the local area are located outside a 0.50-mile radius of the site. Domestic water wells were not identified within 0.25-mile of the site. Therefore, the water supply for the area has not been impacted by Site 333.

### 5.0 CONCLUSIONS AND RECOMMENDATION

The results of B&R Environmental's CA at CSS Site 333 suggest the following:

- Groundwater in the surficial aguifer at the site has a G-II classification;
- Private potable water wells were not identified within 0.25-mile radius of the site;
- Municipal well fields and surface water intakes were not identified within a 0.50-mile radius of the site;
- Analysis of soil samples collected during the Tank Closure Assessment indicate the soils are not hazardous.
- Excessively contaminated soil in vadose soils was not encountered during the CA;
- Oil stain soils identified at the water table surface appear to be from heavily weathered diesel;
- Free product was measured at a thickness of 0.15 foot in well PCY-333-MW04;
- The source of dissolved hydrocarbons appears to be from old spills which occurred at the diesel UST located adjacent to Site 333;
- Analysis of groundwater samples show the benzene, total VOAs, EDB, arsenic, cadmium, chromium, total naphthalene, and 1-2-dichloroethane constituents were reported at concentrations below the State Action Level for No Further Action (NFA) criteria for a G-II aquifer without wells (FDEP, 1990) (Table 5-1). Concentrations of total naphthalene, lead, vinyl chloride, phenanthrene, and cis-1,2-dichloroethane were reported at levels which meet State Action Levels for Monitoring Only status for G-II aquifers without wells.

Based upon the hydrogeological and chemical data presented in this CAR, and the CA criteria for Monitoring Only, the site qualifies for Monitoring Only status. However, we believe prior to developing a Monitoring Only Plan, an Alternate Remedial Procedure should be performed in an attempt to reduce concentration levels of naphthalene, total naphthalene, and TRPH, in source wells TW01 and PCY-333-MW01. The Alternative Remedial Procedure would also address reducing the lead concentration in PCY-333-MW02. An Alternative Remedial Procedure applied to this site may be effective at reducing the constituents in the groundwater to NFA levels or reduce the required monitoring period. A Remedial Action Procedure will be prepared and submitted to the FDEP upon review and approval of this Contamination Assessment Report.

### TABLE 5-1 MAXIMUM ACCEPTABLE GROUNDWATER CONSTITUENT LEVELS **Site 333**

### Coastal Systems Station, Panama City, Florida FDEP Facility No. 038518667

Analyte or Analytical Method	Highest Ground Water Constituent Level in Site Monitoring Wells	No Furt	her Action	Monitoring Only					
			G-II Aquifer	G-II Aquifer wit	G-II Aquifer without wells				
		(with wells)	(without wells)	source	perimeter	source	perimeter		
Total BTEX	15	50	50	500	50	1000	50		
Benzene	2.2	1	50	250	1	500	50		
TRPH	21.5	5^	5^	50^	5^	100^	5^		
Lead	677	50	50	- 500	50	1^	50		
EDB	<0.02	0.02	0.02	0.02	0.02	0.4	0.02		
Total Naphs	670	100	100	1000	100	2000	100		
EPA 610	<4 to<200	DL	DL	10xDL	DL	20xDL	DL		
EPA 601	<1.8	DW-SRLs	DW-SRLs	10xDW-SRLs	DW-SRLs	20xDW-SRLs	DW-SRLs		
Arsenic	4.0	50	50	500	50	1^	50		
Cadmium	<3.3	10	10	100	10	200	10		
Chromium	21.4	50	50	500	50	1^	50		
EPA 624	NCD	DW-SRLs	DW-SRLs	10xDL-SRLs	DW-SRLs	20xDW-SRLs	DW-SRLs		
EPA 625	160	DW-SRLs	DW-SRLs	10xDL-SRLs	DW-SRLs	20xDW-SRLs	DW-SRLs		

#### Notes:

All data in μg/L unless otherwise noted ^ data in mg/L

Source:

Monitoring wells near suspected hydrocarbon source

Perimeter:

Monitoring wells located at perimeter of plume Total Recoverable Petroleum Hydrocarbons

TRPH: Total Naphs:

Sum of naphthalenes and methylnaphthalenes

DW-SRLs:

Drinking Water Standards or Applicable Site Rehabilitation Levels

DL:

**Detection Limit** 

NCD

No Constituents Detected

### 6.0 REFERENCES

ABB Environmental Services, Inc., 1995, <u>RCRA Facility Investigation</u>, <u>Coastal Systems Station</u> Panama City, Florida.

Florida Department of Environmental Protection, October 1990. *No Further Action and Monitoring Only Guidelines for Petroleum Contaminated Sites*, Bureau of Waste Cleanup, Technical Review Section.

Heath Ralph C., 1983, <u>Basic GroundWater Hydrology</u>: US Geological Survey Water Supply Paper 2220.

NAVFAC Drawing No. 504-6560, Oil Spill Control Equipment Clean-Up Wash Rack Plan & Details, August 1, 1978.

Southern Waste Services Environmental First Response, 1996, <u>Closure Assessment Report FAC.</u>
#333 Waste Oil & Oil/Water Separator, Coastal Systems Station, Panama City, Florida.

U.S. Geological Survey, Panama City, FLA., Quadrangle 1982. 7.5 minute series, Topographic Quadrangle Maps of Florida: scale 1:24,000.

U.S. Geological Survey. Panama City Beach, FLA., Quadrangle 1982. 7.5 minute series, Topographic Quadrangle Maps of Florida: scale 1:24,000.

### APPENDIX A

## DISCHARGE NOTIFICATION FORM AND INITIAL REMEDIATION ACTION NOTIFICATION FORM



### DEPARTMENT OF THE NAVY

# COASTAL SYSTEMS STATION DAHLGREN DIVISION NAVAL SURFACE WARFARE CENTER 6703 WEST HIGHWAY 98 PANAMA CITY FL 32407-7001

IN REPLY REFER TO:

5090

Ser 051E/225

Pollutant Storage Tank Program Attn: Mr. Dennis D. Pinkovsky, Ph.D. 619 North Cove Boulevard, Suite C Panama City, FL 32401

Dear Mr. Pinkovsky:

The Discharge Reporting Form provided as enclosure (1) is forwarded as required to report our discovery of contaminated soil during the removal of NSWCCSS waste oil tank #333.

Enclosure (2) is provided for the remedial action we initiated in an attempt to remove all of the excessively contaminated soil. After initiation of the remedial action we concluded from the extent and direction of contamination that the source was not from this tank. It was most likely from an adjacent contaminated site known as Solid Waste Management Unit #1. This adjacent site is an area once used as the NSWCCSS landfill in the late 1940's and early 1950's. It is presently under investigation for possible corrective action under the Navy's Installation Restoration Program. Based on our conclusion we immediately stopped our attempts to remediate the site.

Enclosure (3) is a tank registration update being provided for our removal of this waste oil tank.

If you require additional information at this time, please contact Mr. Mike Clayton, Code 051EMC, at (904) 235-5859 or Mr. Bill Logsdon, Code 051EBL, at (904) 235-5474.

Sincerely,

DY L. GREEN
Assistant Public Works Officer

By dilection of

the Commanding Officer

#### Encl:

- (1) FDEP Form 17-761.900(1)
- (2) Initial Remedial Action Notification
- (3) FDEP form 17-761.900(2)

Copy to:

FDEP (Mr. Eric Nuzie)

NAS Jacksonville (Mr. Jerry Wallmeyer)



### Florida Department of Environmental Regulation

Twin Towers Office Bldg. ● 2600 Blair Stone Road ● Tallahassee, Florida 32399-2400

2 2 1 2 2 2 2 2	
Form Tiss Discharge Reporting Form	_
Effective Date December 10, 1990	-
DER Application No.	_
(Fleed in by DER)	-

### Discharge Reporting Form

Use this form to notify the Department of Environmental Regulation of:

- 1. Results of tank tightness testing that exceed allowable tolerances within ten days of receipt of test result.
- 2. Petroleum discharges exceeding 25 gallons on pervious surfaces as described in Section 17-761.460 F.A.C. within one working day of discovery.
- 3. Hazardous substance (CERCLA regulated), discharges exceeding applicable reportable quantities established in 17-761.460(2) F.A.C., within one working day of the discovery.
- 4. Within one working day of discovery of suspected releases confirmed by: (a) released regulated substances or pollutants discovered in the surrounding area, (b) unusual and unexplained storage system operating conditions, (c) monitoring results from a leak detection method or from a tank closure assessment that indicate a release may have occurred, or (d) manual tank gauging results for tanks of 550 gallons or less, exceeding ten gallons per weekly test or five gallons averaged over four consecutive weekly tests.

Mail to the DER District Office in your area listed on the reverse side of this form

## PLEASE PRINT OR TYPE Complete all applicable blanks

\_\_\_\_\_ 3. Date: 11/28/95 DER Facility ID Number: <u>038518667</u> \_\_\_\_ 2. Tank Number: \_\_\_333 4. Facility Name: Coastal Systems Station Facility Owner or Operator: United States Navv Facility Address: Code 051E, 6703 W Hwy 98, Panama City, FL 32407-7001 Telephone Number: ( 904 ) 235-5859 \_\_\_\_\_ County: BAY Mailing Address: Same as Above 5. Date of receipt of test results or discovery: 11/28/95 \_ month/dav/vear 6. Method of initial discovery. (circle one only) F. Vapor or visible signs of a discharge in the vicinity.

G. Closure: System Removal (explain) D. Emptying and Inspection. A. Liquid detector (automatic or manual) E. Inventory control. B. Vapor detector (automatic or manual) H. Other: C. Tightness test (underground tanks only). 7. Estimated number of gallons discharged: <u>UNKNOWN</u> (E) Unknown 8. What part of storage system has leaked? (circle all that apply) A. Dispenser B. Pipe C. Fitting D. Tank 9. Type of regulated substance discharged. (circle one) A. leaded gasoline D. vehicular diesel (L.) used/waste oil V. hazardous substance includes pesticides, ammonia, chlorine and derivatives (write in name or Chemical Abstract B. unleaded gasoline F. aviation gas M. diesel Service CAS number)\_ C. gasohol G. jet fuel O. new/lube oil Z. other (write in name) \_ 10. Cause of leak. (circle all that apply) I. Other (specify) \_\_\_\_\_ (A) Unknown C. Loose connection E. Puncture G. Spill \_\_ B. Split D. Corrosion F. Installation failure H. Overfill 11. Type of financial responsibility. (circle one) A. Third party insurance provided by the state insurance contractor (C) Not applicable B. Self-insurance pursuant to Chapter 17-769.500 F.A.C. D. None 2. To the best of my knowledge and belief all information submitted on this form jextrue, accurate, and complete. A. OSTER, LCDR. ine or Owner, Operator or Authorized nepresentative

Northwest District 160 Governmental Center Pensecola, :Florida 32501-5794 2014-324 8300

## INITIAL REMEDIAL ACTION NOTIFICATION FORM

This notification provides written confirmation of initial remedial action (IRA) as required by Chapter 17-770.300(5) and (8), Florida Administrative Code. Notification must be within three working days of initiation of an IRA. The notification must be submitted to the appropriate contracted local program and/or:

Florida Department of Environmental Protection
Bureau of Waste Cleanup
Engineering Support Section
2600 Blair Stone Road
Tallahassee, FL 32399-2400
(904) 488-3935

Upon completion of the IRA program task, an Initial Remedial Aciton Report (or its equivalency) should be submitted for technical review.

I.	FACILITY NAME: COASTAL SYSTEMS STATION	
	Facility Address: Code 051E, Coastal Systems Station, 6703 W Hwy 98, Panama Ci	.ty
	DER Facility Number (if applicable): 038518667	
	Date IRA Initiated: 11/28/95 Date IRA Completed: 11/28/95	
ıı.	FREE PRODUCT RECOVERY	:
Α.	Type(s) of Product Discharged: N/A	
В.	Quantity  1. Estimated Gallons Lost: N/A	
c.	Method of Product Recovery: N/A	
D.	Type of Discharge During Product Recovery: N/A	
E.	Type of Treatment, i.e., Oil/Water Separator: and Expected  Effluent Quality from Any Discharge: N/A	
F.	Quantity and Disposal of Recovered Product: N/A	

**MAY 1994** 

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### III. SOIL EXCAVATION

A. Estimated Volume of Excessively Contaminated Soil Excavated in
Cubic Yards: 2 CY
B. Estimated Dimensions of Excavation Including Depth of Excavation(s): 12' long x 9' wide x 5' deep
C. Type(s) of Product in Soil: Waste Petroleum (old)
D. Type of Instrument and Method Used to Determine Excessive
Soil Contamination: Organic Vapor Analyzer (OVA) Instrument with a Flame
Ionization Detector (FID). Samples were screened with and without a carbon filte
•
IV. ADDITIONAL COMMENTS: Initial remediation could not remediate all excessively
contaminated soil due to the location and extent of contamination discovered. The
contamination appears to not have been created by this oily waste management system
but may be that of an adjacent contaminated site being investigated in the Navy's
Installation Restoration Program.
MTCHZET. D. CT.AVTCN
MICHAEL D CLAYTON Print Person Completing Form
Print Person Completing Form
Print Person Completing Form
Print Person Completing Form  Michael D. Clayler 12/01/15 NSWCCSS Public Works Environmental Engineer
Print Person Completing Form    Michael   Class   12/01/15 NSWCCSS Public Works Environmental Engineer Signature, Date Title, Affiliation Code 051E
Print Person Completing Form  Michael Clark 12/01/15 NSWCCSS Public Works Environmental Engineer Signature, Date Title, Affiliation Code 051E 6703 W Hwy 98, Panama City, FL 32407-7001
Print Person Completing Form    Michael   Class   12/01/15 NSWCCSS Public Works Environmental Engineer Signature, Date Title, Affiliation Code 051E
Print Person Completing Form    Michael D. Clay   12/0/15 NSWCCSS Public Works Environmental Engineer Signature Date Title, Affiliation Code 051E 6703 W Hwy 98, Panama City, FL 32407-7001   Company Address
Print Person Completing Form  Michael Clark 12/01/15 NSWCCSS Public Works Environmental Engineer Signature, Date Title, Affiliation Code 051E 6703 W Hwy 98, Panama City, FL 32407-7001



### Florida Depariment of Environmental Regulation

Twin Towers Office Bidg. • 2600 Blair Stone Road • Tallahassee, Florida 32399-2400

Form Tide Storage Tank Registration Form
Effective Date December 10, 1990
DER Application No
(Feed in by DCM)

## Storage Tank Registration Form

		•	
Diagon Drint or Tim	a Daviou Includator	ns Before Completing F	OFm.
Please Print or IVD	e - neview instruction	is before Completing F	ULILL

3. New Re	egistration	New C	wner Dala		ty Revision	] Tank(		X		
Tank(s)	Address: 6	703 West	Hwy 98							
1							-	-	235_585	9
Owner City/Sta	Mailing Addr te/Zip: <u>Par</u>	ess: 6703 nama City	FL 3240	98 7-7001	Station)				235 - 585	9
L	n (optional)	Lalitude: _	Line For E	ach Tank	de:°' At This Fac oks in use; 9	ility (Use	Codes - S		hip	28 / 95 Range
9	10	11	12	13	14	15	16	17	18	19
	550	L .	XX/79	. U	E	В	Х	В	0	11/28/9
*For ne To the best  W. A. OS  Print name	rw tank install of my knowled TER, LCDR & title of own	Certified Contaction or tankedge and be	c removal lief all information WO OFFICED zed person little 17, Sinte 9 200 33	_	ed on this for	Depa	Ecurate and	fessional Reg	2///55 Da 1900 E.G.	ense Number  lle  wheest Oistict ongress Ave., Suite A Besch, Florids 13406 07-432 7850

# APPENDIX B CAR SUMMARY SHEET

_		" AUGESSMERT .	•		·
Fac	ility Name: Coastal Systems Sta		<del>-</del> :		jabursement Site 🔲
	Location: Panana City Fi				_
	EDI #:		8518661	Ot	ner:
	oi/water sopacotor, w				
(1)	Source of spill: diesal took		spill: un ka		
(2)	Type of <u>qasoline group</u>	gasoline lost	<u>ker</u>	osene group	gallons lost
	product: Leaded		← ker	озеле	
	unleaded regular		√ die	sel	unkauwn
	unleaded presiva		=	4 Jet fuel	
	gasohol	·	☐ Jet	A fuel	
		, ,	•		
(3)	Description of IRA (if any): 50:/5 4		☐ Fre	e product resoval	$: \underline{4 \frac{1}{8}} \text{ (gals)}$
	- From work oil took pit		$\square$	Soil removal	: (cubic yds
				Soil incineration	: (cubic yds
(4)	Free product still present? (yes/no)	Kaximum appari	ent product this	kness; <u>0./5</u>	_ (ft)
					EDB: 40.02
	contaminant levels (ppb):				other: TRPH
	/ Te	tal Manh Hanlows:	670 (14	10-1111	
(6)	Brief lithologic description:	in yi chloride: 1.4	phenant	hrene:160	
(8) (9)	Areal and vertical extent of soils contains the current soil concentration (OVA:	Depth of verti Silv Sompled er sampling: 11/25	or (EPA Method cal contamination 7////2	on: <u>Vertical Exten</u>	t will not instilled
	QAPP approved? (yes/no) Date: 6//6	•	6 11. d. 18:0	···· 3-1	2- V )
	Direction (e.g. NNW) of surficial ground		- XAPAGWST (Fig	ure <u>3 1</u> on	page
	Average depth of groundwater: 4.5 4.		- (so	100 000	\
			<i>(13</i>	tweet peg.	1-13)
	Estimated rate of groundwater Box. 1.18  Hydraulic gradient across site:				
			linda.	Manha	
(10)	Aquifer characteristics:  Hydraulic conductivity	<u>Value</u>	Units It day	Method Share does	<del></del> , ,
	Storage coefficient	<u> 7.L</u>	117 974 	Slug test	CHUN
	Aquifer thickness		7+	Begin to w	ter subfracted fro
	Effective soil porosity	•30	<del></del>	Literatury	10/10
	Transmissivity	156	7+2/day	T= Kb	0-700
			——————————————————————————————————————	٠	/ 1 1
17)	other remarks: Solid Waste Monage				
	upgradient of	woste oil	tonk Dies	sel tonk soo	ce for
	old product observ	red at mater a	table/ Soil	interpres	

# APPENDIX C TANK CLOSURE ASSESSMENT



### DEPARTMENT OF THE NAVY

COASTAL SYSTEMS STATION DAHLGREN DIVISION
NAVAL SURFACE WARFARE CENTER
6703 WEST HIGHWAY 98
PANAMA CITY FL 32407-7001

IN REPLY REFER TO:

5090 Ser 051E/027

0 6 MAR 1996

Pollutant Storage Tank Program Attn: Mr. Dennis D. Pinkovsky, Ph.D. HRS Environmental Health Services 619 North Cove Boulevard, Suite C Panama City, FL 32401

Dear Mr. Pinkovsky:

Enclosure (1) is provided regarding the excavation and removal of the underground waste oil tank #333 and the oil/water separator associated with the Coastal Systems Station (NSWCCSS) Spill Containment Boom wash area, facility #333.

Petroleum contaminated soils discovered during the excavation were verified by using an Organic Vapor Analysis. Initial remediation began but soon stopped as it was evident the amount of remediation needed would greatly exceed the scope of work identified in the tank removal contract. The initial remedial action report will soon follow once disposal of the contaminated soil has been completed.

NSWCCSS will be using the services of Brown and Root Environmental, a Division of Brown & Root, Inc., through a contract with the Naval Facilities Engineering Command, Southern Division, to conduct the petroleum contamination assessment. This contamination assessment will be conducted in accordance with the requirements of the Florida Petroleum Contamination Agreement signed in October 1990 by the Secretary of the Florida Department of Environmental Regulation and the Assistant Secretary of the Navy.

5090 Ser 051E/027

Should you have any questions concerning the enclosure please contact Mr. Mike Clayton, Code 051EMC, at (904) 235-5859 or Mr. Bill Logsdon, Code 051EBL, at (904) 235-5474.

Sincerely,

W. A. OSTER

Lieutenant Commander, U.S. Navy

By direction of

the Commanding Officer

(1) Closure Assessment Report, Facility #333

Copy to:

FDEP

SOUTHDIV (Mr. Nick Ugolini, Code 1843) Brown & Root Environmental (Mr. Gerald Goode)



CLOSURE ASSESSMENT REPORT
FAC. #333 WASTE OIL TANK & OIL/WATER SEPARATOR
COASTAL SYSTEMS STATION
PANAMA CITY, FLORIDA

Prepared By:

SWS Environmental First Response January 2, 1996

ENCL (1)

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agg.A	endix 1	B - CHAIN-OF CUSTODY, LAB ANALYSES	

#### 1.0 INTRODUCTION

This is a report summarizing SWS's work in completing an assessment of the soils related to the removal of a 550 gallon underground storage tank (UST), oil water separator, and related underground piping at the Naval Surface Warfare Center Coastal Systems located in Panama City Beach, Florida. The work was completed at the site adjacent to the concrete wall labeled area #333, November 28, 1995 (see site map attached).

#### 1.1 SCOPE OF WORK

SWS's investigative efforts included the following elements within the scope of work:

- 1. Excavated Area Assessment and organic vapor analysis (OVA) of soils.
- 2. Sampling and Lab Analyses of: a) a composite of waste oil tank Soils b) soils located three feet east of the waste oil tank c) soils adjacent to the oil water separator location and d) a composite soil sample of the excavated soils.
- 3. Sampling of groundwater if contaminants are found present in the soil during UST removal.
- 4. Completion of this report.

### 2.0 EXCAVATED AREA ASSESSMENT

As mentioned above, Southern Waste Services was mobilized to the naval facility to assess the contaminated soils, and transport them to a proper disposal facility. Upon arrival at the site ground water was noticed to be at 5.5 feet below land surface(BLS)at the bottom of the excavation and above a concrete slab upon which the underground storage tank had rested (See Appendix A, Field Notes). All associated piping along with the UST and oil/water separator had been subsequently removed and excavated soil piled on top of visquene and covered to ensure no migration due to precipitation.

### 2.1 Soil Sampling Within the UST Pit

As shown on the site map (Attached) four sampling points labeled S-1, S-2, S-3, S-4, were picked at the corners of the UST excavation at approximately four feet BLS (See Site Map, Attached). All

samples showed corrected values less than 10 ppm excessive hydrocarbon contamination in soil except for sample S-3 which exhibited a corrected value of 950 ppm. The soil sample at S-3 appeared to have a greenish dark gray oily sheen. Further investigation of the contaminated soil zone indicated that it may be within a groundwater - vadose "smear" zone.

2.2 Soil Sampling within the Pipeline trench and Oil/Water Separator

Three soil sampling points were located along the pipeline trench and within the oil/water separator area. As Shown on the site map they all exhibited a corrected value of less than 10 ppm excessive hydrocarbon contamination in the soil.

2.3 Exploratory Boring of Area between Containment and UST Area

Hand auger borings #8 and #9 were installed to the top of ground water found to be approximately five feet BLS. Both Hand auger borings found clean soil to the depth of 4.5 feet BLS. Hand auger boring #8 located 3.5 feet east of sample S-3 discovered contaminated soil at 4.5 feet BLS with a corrected organic reading of 2100 ppm (See Field Notes, Appendix A). Hand auger boring #9 located adjacent to the fenced in waste and drum containment area found greenish, gray oily soil 4.5 feet BLS. Soil samples taken from the contaminated zone gave readings of 900 ppm for total hydrocarbons and 1200 ppm filtered hydrocarbon for a corrected value of -300 ppm. This is indicative of a break through in the All soil samples appeared to dramatically increase in their corrected hydrocarbon levels at 4.5 feet BLS, diesel or old hydrocarbon fuel odors were very apparent. Soil samples from the 4.5 foot deep zone were taken for lab analysis from Hand auger boring #9 (see the following lab analysis discussion).

### 3.0 SAMPLING AND LAB ANALYSES OF SITE SOILS

During the course of the field sampling, soil samples for laboratory analysis were collected from soils within the UST excavated area adjacent to sample sites S-1 through S-4. All samples of environmental media were collected in accordance with SWS's State approved CompQAPP #920203. All samples were properly contained, labeled and placed on ice for transport to the laboratory under chain-of-custody (See Appendix B). Results of analyses are discussed below.

### 3.1 Soil Sampling Within the UST (Waste Oil Tank) Pit

A composite soil sample for lab analysis was taken from borings S-1, S-2, S-3 and S-4 from a depth of four feet BLS and composited (see Site Map, Appendix B, Chain of Custody, Sample 001). These samples were analyzed by GEOS, Inc. for *Volatile Organics* (EPA Method 8240), *Semivolatile Organics* (EPA Method 8270), *Total Recoverable Petroleum Hydrocarbons* (EPA Method 9073) and *RCRA Metals*.

### 3.2 Soil Sampling adjacent to oil/water Separator Area

A soil sample S-004 was taken from the excavation wall adjacent to the oil/water separator by field sample location S-7 and subjected to lab analyses mentioned above (See Site Map attached, sample S-004).

### 3.3 Soil Sampling of Unscheduled Location

A third unscheduled sample S-002 (see Site Map, Appendix B, Lab Analyses) was taken for lab analyses EPA Methods 8240, 8270, 9073 and RCRA Metals from Hand auger boring S-9 and the potentially contaminated soil zone at 4.5 feet BLS.

### 3.4 Composite soil Sample of Excavated Soil

A composite soil sample S-003 was taken from the excavated soils for lab analysis including TCLP Volatiles and TCLP Metals for disposal characterization.

### 4.0 LAB ANALYSES OF SOIL SAMPLES

Analyses from the composite soil sample S-001 taken at the waste oil tank area exhibited abnormal concentrations of RCRA metals, lead 30.4 mg/Kg, barium 2.3 mg/Kg, and chromium 2.7 mg/Kg. All analyses for organic constituents this sample returned below detection limits (BDL), except for total recoverable hydrocarbons (TRPH) 960 mg/Kg, 1-methyl naphthalene 4800  $\mu g/Kg$ , 2-methyl naphthalene 7500  $\mu g/Kg$ , and naphthalene 1300  $\mu g/Kg$  for total naphthalene 13,600  $\mu g/Kg$  or 13.6 mg/Kg(see lab analyses Appendix B).

Analyses from the soil sample S-002 taken at Hand auger boring S-9 (See attached Site Map) located three feet east of the waste oil tank area exhibited concentrations of RCRA metals, barium 1.4 mg/Kg, and chromium 3.2 mg/Kg. All analyses for organic constituents this sample returned BDL except for TRPH 320 mg/Kg, 1-methyl naphthalene 520 µg/Kg, and 2-methyl naphthalene 410 µg/Kg for total naphthalenes 930 µg/Kg (see lab analyses Appendix B).

Analyses from the soil sample S-004 taken beside soil boring S-7 (See attached Site Map) located adjacent to the oil/water separator excavation area exhibited concentrations of RCRA metals, barium 26 mg/Kg, cadmium 1.0 mg/Kg, chromium 3.2 mg/Kg, silver 3.8 mg/Kg, and mercury 0.104 mg/Kg. All analyses for organic constituents this sample returned below detection limits except for TRPH 12 mg/Kg (see lab analyses Appendix B).

### 5.0 PLACEMENT AND TESTING OF INITIAL MONITOR WELL

Monitor well MW-1 was placed December 12, 1995. The location is shown on the site map attached. MW-1 was placed to test conditions in the surficial aquifer. The temporary monitor was screened from three feet BLS to eight feet BLS, extending both above and below the water table found at approximately 4.5 feet BLS.

Following the placement of temporary monitor well MW-1, a ground water sample was collected December 15, 1995. Lab testing data and Chain-of-Custody for this sampling date are included in Appendix B. The levels of contamination that were detected during this initial sampling were found to be high for benzene 1.5  $\mu$ g/L. The "Monitor Only" for Perimeter well in G-II groundwater criteria for benzene is 1  $\mu$ g/L if public or private drinking water wells are located within a half mile or quarter mile radius respectively.

#### 6.0 CONCLUSION

Under the limits of the specified scope of work listed in the proposal and contract for this environmental cleanup, the site has shown indication of contamination by petroleum products at the UST area and adjacent Hand auger boreholes S-8 & S-9. The constituents discovered at this location are believed to have resulted from contamination migrating within the "smear zone caused by rise and fall of the surficial aguifer.

This concludes SWS's report of the environmental assessment and cleanup at Area 333, Naval Surface Warfare Center Coastal Systems Station, Panama City Beach, Florida.

### 7.0 CERTIFICATION

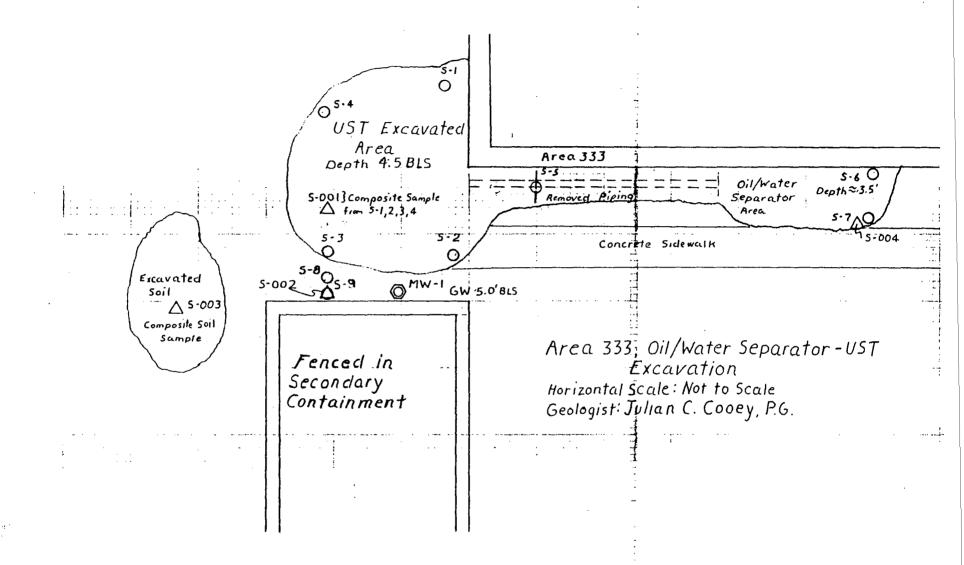
This Contamination Assessment Report was prepared by or under the personal direction of the undersigned registered professional. All parts of this Plan that are concerned with the practice of professional geology were prepared by Mr. Julian C. Cooey, P.G. Field sampling was conducted under the Florida State approved comprehensive quality assurance project plan (CompQAPP #920203). The site-specific health and safety plans were prepared by Mr. Julian C. Cooey, a certified Site Safety Supervisor and certified Environmental Trainer, per 29 CFR 1910.

Respectfully Submitted,

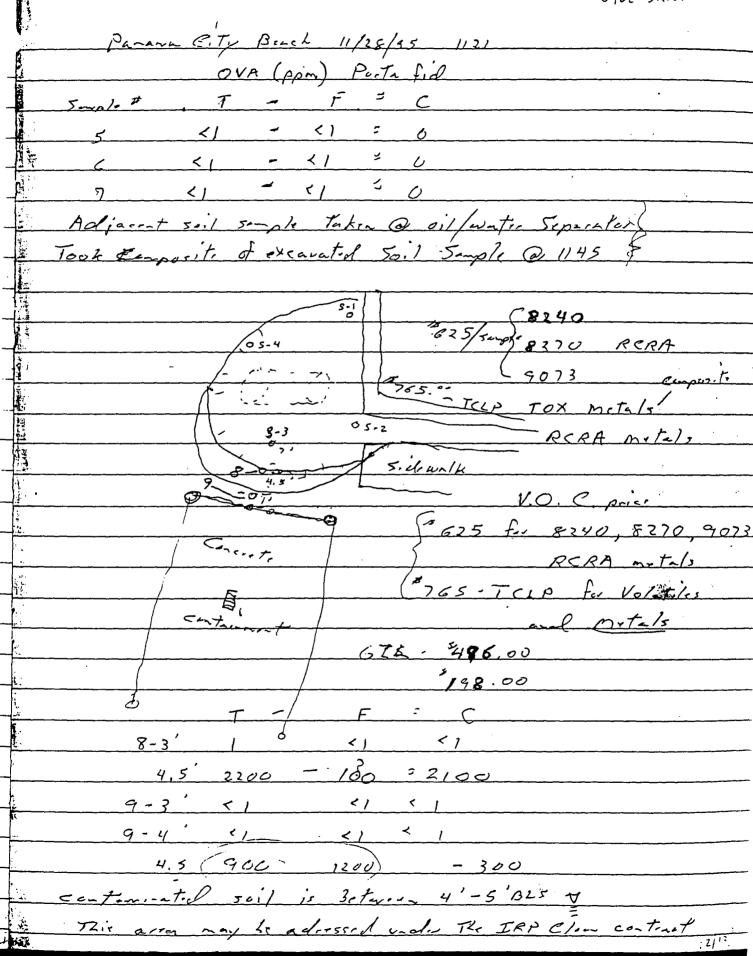
Julian C. Cooey, P.G., CET Florida Registration #32

ATTACHMENT - 1 SITE MAP

Constant



APPENDIX A FIELD NOTES



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APPENDIX B
CHAIN-OF CUSTODY
LAB ANALYSES

- 1				
	CHAIN	OF	CUSTODY	RECORD
	-,			

G	A	N	2	Inc
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1057 N. ELLIS ROAD, SUITE 17, JACKSONVILLE, FL 32254-2249 • (904) 786-8340 [] 5909A BRECKENRIDGE PARKWAY, TAMPA, FL 33610-4237 • (813) 626-0101

1	JOB NUMBER	1/2
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CORPORATE OFFICES
1057 N. ELLIS ROAD, SUITE 17
JACKSONVILLE, FL 32254-2249

(904) 786-8340 (800) 770-4367 (GEOS) FAX: (904) 786-7489

OLOGICAL, ENVIRONMENTAL AND OCEANOGRAPHIC SCIENCES

ANALYTICAL LABORATORY 1627 EAST 8th STREET JACKSONVILLE, FL 32206

(904) 354-6755 FAX: (904) 354-3799

S0U02

Attn: HARRY MARSH

SOUTHERN WASTE SERVICES, INC. 1619 MOYLAN ROAD PANAMA CITY BEACH, FL. 32407 (800)852-8878 Page 1 6 Dec 1995

Report J5-11-080-01 LAB ID. 82223/E82101

Sample Description:

NCSC/JLS CONTRACTOR

PANAMA CITY BEACH, FL.

WASTE OIL TANK COMPOSITE SOIL SAMPLE

P.O. #: HM 95-2570

SAMPLE ID.: SAMPLE 001 COLLECTED: 11/28/95 09:50

RECEIVED: 11/28/95 09:50

COLLECTED BY: J. COOEY

Parameter	Result	Units	Method	Det. Limit	Extracted	Analyzed	Analyst
Hydrocarbons, Total IR	960	mg/tg	9073	0.2	12/04/95	12/05/95	AM
ICAP Metals, Total			3050/6010				
Arsenic	BOL	mg/kg		7.5	11/30/95	11/30/95	1C
Barium	2.3	mg/kg		0.3	11/30/95	11/30/95	JC
Cadmium	BDL	mg/kg		0.5	11/30/95	11/30/95	JC
Chromium	2.7	mg/kg		0.5	11/30/95	11/30/95	JC
Lead	30.4	mg/kg		10.0	11/30/95	11/30/95	JC
Selenium	BDL	mg/kg		10.0	11/30/95	11/30/95	1C
Silver	BDL	mg/kg		1.0	11/30/95	11/30/95	1C
Mercury, Total	<0.075	mg/kg	7471	0.075	12/01/95	12/01/95	JC
Semi-Volatile Organics			3550\8270	-	•		
Acenaphthene	BDL	μg/Kg		330	12/04/95	12/04/95	<b>A</b> T
Acenaphthylene	BDL	μg/Kg		330	12/04/95	12/04/95	ΑT
Aldrin	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Aniline	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Anthracene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Benzidine	BDL	μg/Kg		660	12/04/95	12/04/95	AT
Benzoic acid	BDL	μg/Kg		1700	12/04/95	12/04/95	AT
Benzo (a) anthracene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Benzo (b) fluoranthene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Benzo (k) fluoranthene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Benzo (g,h,i) perylene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Benzo (a) pyrene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Benzyl alcohol	BDL	μg/Kg		330	12/04/95	12/04/95	AT
а-ВНС	BDL	μg/Kg		330	12/04/95	12/04/95	AT
b-BHC	BDL	μg/Kg		330	12/04/95	12/04/95	AT
d-BHC	BDL	μg/Kg		330	12/04/95	12/04/95	AT
g-BHC (Lindane)	BDL	µs/Kg		330	12/04/95	12/04/95	AT
Bis (2-chloroethoxy) methane	<b>3</b> 0 L	, #9/Kg		330	12/04/95	12/04/95	AT
Bis (2-chloroethy() ether	301	μg/Xg		330	12/04/95	12/04/95	AT.
Bis (2-chloroisopropyl) ether	BDL	μg/Kg		330	12/04/95	12/04/95	AT

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arameter	Result	Units	Method	Det. Limit	Extracted	Analyzed	Analyst
Bis (2-ethylhexyl) phthalate	BDL	μg/Kg		330	12/04/95	12/04/95	AT
4-Bromophenyl phenyl ether	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Butylbenzylphthalate	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Chlordane	BDL	μg/Kg		1700	12/04/95	12/04/95	AT
4-Chloroaniline	BDL	μg/Kg		330	12/04/95	12/04/95	AT
4-Chloro-3-methylphenol	BDL	μg/Kg		330	12/04/95	12/04/95	AT
2-Chloronaphthalene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
2-Chlorophenol	BDL	μg/Kg		330	12/04/95	12/04/95	AT
4-Chlorophenyl phenyl ether	BDL	μg/Kg		<b>3</b> 30	12/04/95	12/04/95	AT
Chrysene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
4,4'-DDD	BDL	μg/Kg		330	12/04/95	12/04/95	AT
4,4'-DDE	BDL	μg/Kg		330	12/04/95	12/04/95	AT
4,4'-DDT	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Dibenzo (a,h) anthracene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Dibenzofuran	BDL	μg/Kg		330	12/04/95	12/04/95	ΑT
Di-n-butylphthalate	BDL	μg/Kg		330	12/04/95	12/04/95	AT
1,2-Dichlorobenzene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
1,3-Dichlorobenzene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
1,4-Dichlorobenzene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
3,3'-Dichlorobenzidine	BDL	μg/Kg		660	12/04/95	12/04/95	AT
2,4-Dichlorophenol	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Dieldrin	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Diethylphthalate	BDL	μg/Kg		330	12/04/95	12/04/95	AT
2,4-Dimethylphenol	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Dimethylphthalate	BDL	μg/Kg		330	12/04/95	12/04/95	AT .
2,4-Dinitrophenol	BDL	μg/Kg		1700	12/04/95	12/04/95	AT
2,4-Dinitrotoluene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
2,6-Dinitrotoluene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Di-n-octylphthalate	BOL	μg/Kg		330	12/04/95	12/04/95	AT
1,2-Diphenylhydrazine	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Endosulfan I	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Endosulfan II	80L	μg/Kg		330	12/04/95	12/04/95	AT
Endosulfan sulfate	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Endrin	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Endrin aldehyde	BDL	μg/Kg		330	12/04/95	12/04/95	AT

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SOUTHERN WASTE SERVICES, INC. 1619 MOYLAN ROAD PANAMA CITY BEACH, FL. 32407 (800)852-8878 Page 3 6 Dec 1995 Report J5-11-080-01 LAB ID. 82223/E82101

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arameter	Result	Units	Method	Det. Limit	Extracted	Analyzed .	Analyst
Fluoranthene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Fluorene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Heptachlor	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Heptachlor epoxide -	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Hexach lorobenzene	BDL.	μg/Kg		330	12/04/95	12/04/95	AT
Hexachlorobutadiene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Hexachlorocyclopentadiene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Hexachloroethane	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Indeno (1,2,3-c,d) pyrene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Isophorone	BDL	μg/Kg		330	12/04/95	12/04/95	AT
2-Methyl-4,6-dinitrophenol	BDL	μg/Kg		1700	12/04/95	12/04/95	AT
1-Methylnaphthalene	4800	μg/Kg		330	12/04/95	12/04/95	AT
2-Methylnaphthalene	7300	μg/Kg		330	12/04/95	12/04/95	AT
2-Methylphenol	BDL	μg/Kg		330	12/04/95	12/04/95	AT
4-Methylphenol	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Naphthalene	1300	μg/Kg		330	12/04/95	12/04/95	AT
2-Nitroaniline	BDL	μg/Kg		1700	12/04/95	12/04/95	AT
3-Nitroaniline	BDL	μg/Kg		1700	12/04/95	12/04/95	AT
4-Nitroaniline	BDL	μg/Kg		1700	12/04/95	12/04/95	AT
Nitrobenzene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
2-Nitrophenol	BOL	μg/Kg		330	12/04/95	12/04/95	AT
4-Nitrophenol	BDL	μg/Kg		1700	12/04/95	12/04/95	AT
N-Nitrosodimethylamine	BDL	μg/Kg		330	12/04/95	12/04/95	AT
N-Nitrosodiphenylamine	BDL	μg/Kg		330	12/04/95	12/04/95	AT
N-Nitrosodi-n-propylamine	BDL	μg/Kg		330	12/04/95	12/04/95	AT ·
PCB-1016	BDL	μg/Kg		1700	12/04/95	12/04/95	AT
PCB-1221	BDL	μg/Kg		1700	12/04/95	12/04/95	AT
PCB-1232	BOL	μg/Kg		1700	12/04/95	12/04/95	AT
PCB-1242	BDL	μg/Kg		1700	12/04/95	12/04/95	AT
PCB-1248	BDL	μg/Kg		1700	12/04/95	12/04/95	AT
PCB-1254	BDL	μg/Kg		1700	12/04/95	12/04/95	AT
PCB-1260	BDL.	μg/Kg		1700	12/04/95	12/04/95	AT
Pentachlorophenol	BDL	. μg/Kg		1700	12/04/95	12/04/95	AT
Phenanthrene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Phenol	BDL	μg/Kg		330	12/04/95	12/04/95	AT

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Parameter	Result		Units	Method	Det. Limit	Extracted	Analyzed /	Analyst
Pyrene	BDL		μg/Kg		330	12/04/95	12/04/95	AT
Pyridine	BDL		μg/Kg		330	12/04/95	12/04/95	AT
Toxaphene	BDL		μg/Kg		3300	12/04/95	12/04/95	AT
1,2,4-Trichlorobenzene -	BDL		μg/Kg		330	12/04/95	12/04/95	AT
2,4,5-Trichlorophenol	BDL		μg/Kg		330	12/04/95	12/04/95	AT
2,4,6-Trichlorophenol	BDL		μg/Kg		330	12/04/95	12/04/95	AT
Surrogates								
Nitrobenzene-d5	53	Min: 23		Max:	120			
2-Fluorobiphenyl	77	Min: 30		Max:	115			
Terphenyl-d14	65	Min: 18		Max:	137			
Phenol-d5	51	Min: 24		Max:	113			
2-Fluorophenol	65	Min: 25		Max:	121			
2,4,6-Tribromophenol	69	Min: 19		Max:	122			
/olatile Organics				8240\8260	-			
Acetone	BDL		μg/Kg		250	11/30/95	11/30/95	AT
Benzene	BDL		μg/Kg		10	11/30/95	11/30/95	AT
Bromodichloromethane	BDL		μg/Kg		10	11/30/95	11/30/95	AT
Bromoform	BDL		μg/Kg		10	11/30/95	11/30/95	AT
Bromomethane	BDL		μg/Kg		10	11/30/95	11/30/95	AT .
2-Butanone	BDL		μg/Kg		250	11/30/95	11/30/95	AT
Carbon disulfide	BDL		μg/Kg		10	11/30/95	11/30/95	AT
Carbon tetrachloride	BDL		μg/Kg		10	11/30/95	11/30/95	AT
Chlorobenzene	BDL		μg/Kg		10	11/30/95	11/30/95	AT
Chlorodibromomethane	BDL		μg/Kg		10	11/30/95	11/30/95	AT
Chloroethane	BDL		μg/Kg		10	11/30/95	11/30/95	AT ·
Chloroform	BDL		μg/Kg		10	11/30/95	11/30/95	AT
Chloromethane	BDL		μg/Kg		10	11/30/95	11/30/95	AT
Dibromomethane	BDL		μg/Kg		10	11/30/95	11/30/95	AT
1,4-Dichloro-2-butene	BDL		μg/Kg		10	11/30/95	11/30/95	AT
Dichlorodifluoromethane	BDL		μg/Kg		10	11/30/95	11/30/95	AT
1,1-Dichloroethane	BDL		μg/Kg		10	11/30/95	11/30/95	AT
1,2-Dichloroethane	BDL		μg/Kg		10	11/30/95	11/30/95	AT
1,1-Dichloroethene	BDL		μg/Kg		10	11/30/95	11/30/95	AT
trans-1,2-Dichloroethene	BDL		μg/Kg		10	11/30/95	11/30/95	AT
1,2-Dichloropropane	BDL		μg/Kg		10	11/30/95	11/30/95	AT

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Attn: HARRY MARSH

SOUTHERN WASTE SERVICES, INC. 1619 MOYLAN ROAD PANAMA CITY BEACH, FL. 32407 (800)852-8878 Page 5 6 Dec 1995 Report J5-11-080-01 LAB ID. 82223/E82101

Parameter	Result	Units	Method	Det. Limit	Extracted	Analyzed	Analyst
cis-1,3-Dichloropropene	BDL	μg/Kg		10	11/30/95	11/30/95	AT
trans-1,3-Dichloropropene	BDL	μg/Kg		10	11/30/95	11/30/95	AT
Ethanol	BDL	μg/Kg		1000	11/30/95	11/30/95	AT
Ethylbenzene -	BDL	μg/Kg		10	11/30/95	11/30/95	AT
Ethyl methacrylate	BDL	μg/Kg		10	11/30/95	11/30/95	AT
2-Hexanone	BDL	μg/Kg		100	11/30/95	11/30/95	AT
Iodomethane	BDL	μg/Kg		10	11/30/95	11/30/95	AT
Methylene chloride	BDL	μg/Kg		10	11/30/95	11/30/95	AT
4-Methyl-2-pentanone	BDL	μg/Kg		100	11/30/95	11/30/95	AT
Styrene	BDL	μg/Kg		10	11/30/95	11/30/95	AT
1,1,2,2-Tetrachloroethane	BDL	μg/Kg		10	11/30/95	11/30/95	AT
Tetrachloroethene	BDL	μg/Kg		10	11/30/95	11/30/95	AT
Toluene	BDL	μg/Kg		10	11/30/95	11/30/95	AT
1,1,1-Trichloroethane	BDL	μg/Kg		10_	11/30/95	11/30/95	AT
1,1,2-Trichloroethane	BDL	μg/Kg		10	11/30/95	11/30/95	AT
Trichloroethene	BDL	μg/Kg		10	11/30/95	11/30/95	AT
Trichlorofluoromethane	BDL	μg/Kg		10	11/30/95	11/30/95	AT
1,2,3-Trichloropropane	BDL	μg/Kg		10	11/30/95	11/30/95	AT
Vinyl chloride	BDL	μg/Kg		10	11/30/95	11/30/95	AT
Xylenes (Total)	BDL	μg/Kg		10	11/30/95	11/30/95	AT
Surrogates							
Toluene-d8	101	Min: 81	Max: 1	17			
4-Bromofluorobenzene	90	Min: 74	Max: 12	21			
1,2-Dichloroethane-d4	96	Min: 70	Max: 12	21			

Greg D. Johnson, Lab Director

SOU02

Attn: HARRY MARSH

SOUTHERN WASTE SERVICES, INC. 1619 MOYLAN ROAD PANAMA CITY BEACH, FL. 32407 (800)852-8878 Page 6 6 Dec 1995

Report J5-11-080-02 LAB ID. 82223/E82101

Sample Description:

NCSC/JLS CONTRACTOR PANAMA CITY BEACH, FL.

SOIL SAMPLE 4.5' BLS + 3'E OF W.O.T.

P.O. #: HM 95-2570

SAMPLE ID.: SAMPLE 002 COLLECTED: 11/28/95 11:00

RECEIVED: 11/28/95 COLLECTED BY: J. COOEY

Parameter	Result	Units	Method	Det. Limit	Extracted	Analyzed	Analyst
Hydrocarbons, Total IR	320	mg/kg	9073	0.2	12/04/95	12/05/95	АМ
ICAP Metals, Total			3050/6010				
Arsenic	BDL	mg/kg		7.5	11/30/95	11/30/95	JC
Barium	1.4	mg/kg		0.3	11/30/95	11/30/95	JC
Cadmīum	BDL	mg/kg		0.5	11/30/95	11/30/95	JC
Chromium	3.2	mg/kg		0.5	11/30/95	11/30/95	JC
Lead	BDL	mg/kg		10.0	11/30/95	11/30/95	1C
Selenium	BDL	mg/kg		10.0	11/30/95	11/30/95	JC
Silver	BDL	mg/kg		1.0	11/30/95	11/30/95	JC
Mercury, Total	<0.075	mg/kg	7471	0.075	12/01/95	12/01/95	JC
Semi-Volatile Organics			3550\8270	•			
Acenaphthene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Acenaphthylene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Aldrin	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Aniline	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Anthracene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Benzidine	BDL	μg/Kg		660	12/04/95	12/04/95	AT
Benzoic acid	BDL	μg/Kg		1700	12/04/95	12/04/95	AT
Benzo (a) anthracene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Benzo (b) fluoranthene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Benzo (k) fluoranthene	BDL	μg/Kg		330	12/04/95	12/04/95	AT .
Benzo (g,h,i) perylene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Benzo (a) pyrene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Benzyl alcohol	BDL	μg/Kg		330	12/04/95	12/04/95	AT
a-BHC	BDL	μg/Kg	•	330	12/04/95	12/04/95	AT
b-BHC	BDL	μg/Kg		330	12/04/95	12/04/95	AT
d-BHC	BDL	μg/Kg		330	12/04/95	12/04/95	AT
g-BHC (Lindane)	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Bis (2-chloroethoxy) methane	BDL	. μg/Kg		330	12/04/95	12/04/95	AT
Bis (2-chloroethyl) ether	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Bis (2-chloroisopropyl) ether	BDL	μg/Kg		330	12/04/95	12/04/95	AT

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SOUTHERN WASTE SERVICES, INC. 1619 MOYLAN ROAD PANAMA CITY BEACH, FL. 32407 (800)852-8878 Page 7 6 Dec 1995 Report J5-11-080-02 LAB ID. 82223/E82101

arameter	Result	Units	Method	Det. Limit	Extracted	Analyzed	Analyst
Bis (2-ethylhexyl) phthalate	BDL	μg/Kg		330	12/04/95	12/04/95	AT
4-Bromophenyl phenyl ether	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Butylbenzylphthalate	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Chlordane -	BDL	μg/Kg		1700	12/04/95	12/04/95	AT
4-Chloroaniline	BDL	μg/Kg		330	12/04/95	12/04/95	AT
4-Chloro-3-methylphenol	BDL	μg/Kg		330	12/04/95	12/04/95	AT .
2-Chloronaphthalene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
2-Chlorophenol	BDL	μg/Kg		330	12/04/95	12/04/95	AT
4-Chlorophenyl phenyl ether	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Chrysene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
4,4'-DDD	BDL	μg/Kg		330	12/04/95	12/04/95	AT
4,4'-DDE	BDL	μg/Kg		330	12/04/95	12/04/95	AT
4,4'-DDT	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Dibenzo (a,h) anthracene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Dibenzofuran	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Di-n-butylphthalate	BDL	μg/Kg		330	12/04/95	12/04/95	AT
1,2-Dichlorobenzene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
1,3-Dichlorobenzene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
1,4-Dichlorobenzene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
3,3'-Dichlorobenzidine	BDL	μg/Kg		660	12/04/95	12/04/95	AT
2,4-Dichlorophenol	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Dieldrin	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Diethylphthalate	BDL	μg/Kg		330	12/04/95	12/04/95	AT
2,4-Dimethylphenol	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Dimethylphthalate	BDL	μg/Kg		330	12/04/95	12/04/95	AT
2,4-Dinitrophenol	BDL	μg/Kg		1700	12/04/95	12/04/95	AT
2,4-Dinitrotoluene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
2,6-Dinitrotoluene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Di-n-octylphthalate	BDL	μg/Kg		330	12/04/95	12/04/95	AT
1,2-Diphenylhydrazine	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Endosulfan I	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Endosulfan II	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Endosulfan sulfate	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Endrin	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Endrin aldehyde	BDL	μg/Kg		330	12/04/95	12/04/95	AT

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SOUTHERN WASTE SERVICES, INC. 1619 MOYLAN ROAD PANAMA CITY BEACH, FL. 32407 (800)852-8878 Page 8 6 Dec 1995 Report J5-11-080-02 LAB ID. 82223/E82101

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arameter	Result	Units	Method	Det. Limit	Extracted	Analyzed	Analyst
Fluoranthene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Fluorene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Heptachlor	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Heptachlor epoxide -	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Hexach l orobenzene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Hexachlorobutadiene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Hexachlorocyclopentadiene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Hexachloroethane	BDL	μg/Kg		330	12/04/95	12/04/95	- AT
Indeno (1,2,3-c,d) pyrene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Isophorone	BDL	μg/Kg		330	12/04/95	12/04/95	AT
2-Methyl-4,6-dinitrophenol	BDL	μg/Kg		1700	12/04/95	12/04/95	AT
1-Methylnaphthalene	520	μg/Kg		330	12/04/95	12/04/95	AT
2-Methylnaphthalene	410	μg/Kg		330	12/04/95	12/04/95	AT
2-Methylphenol	BDL	μg/Kg		330	12/04/95	12/04/95	AT
4-Methylphenol	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Naphthalene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
2-Nitroaniline	BDL	μg/Kg		1700	12/04/95	12/04/95	AT
3-Nitroaniline	BDL .	μg/Kg		1700	12/04/95	12/04/95	AT
4-Nitroaniline	BDL	μg/Kg		1700	12/04/95	12/04/95	AT
Nitrobenzene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
2-Nitrophenol	BDL	μg/Kg		330	12/04/95	12/04/95	AT
4-Nitrophenol	BDL	μg/Kg		1700	12/04/95	12/04/95	AT
N-Nitrosodimethylamine	BDL	μg/Kg		330	12/04/95	12/04/95	AT
N-Nitrosodiphenylamine	BDL	μg/Kg		330	12/04/95	12/04/95	AT
N-Nitrosodi-n-propylamine	BDL	μg/Kg		330	12/04/95	12/04/95	AT .
PCB-1016	BDL	μg/Kg		1700	12/04/95	12/04/95	AT
PCB-1221	BDL	μg/Kg		1700	12/04/95	12/04/95	AT
PCB-1232	BDL	μg/Kg		1700	12/04/95	12/04/95	ΑT
PCB-1242	BDL	μg/Kg		1700	12/04/95	12/04/95	AT
PCB-1248	BDL	μg/Kg		1700	12/04/95	12/04/95	AT
PCB-1254	BDL	μg/Kg		1700	12/04/95	12/04/95	AT
PCB-1260	BDL	μg/Kg		1700	12/04/95	12/04/95	AT
Pentachlorophenol	BDL	. μg/Kg		1700	12/04/95	12/04/95	AT
Phenanthrene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Phenol	BDL	μg/Kg		330	12/04/95	12/04/95	AT

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SOUTHERN WASTE SERVICES, INC. 1619 MOYLAN ROAD PANAMA CITY BEACH, FL. 32407 (800)852-8878 Page 9 6 Dec 1995 Report J5-11-080-02 LAB ID. 82223/E82101

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Parameter	Result	U	ni ts	Method	Det.	Limit	Extracted	Analyzed	Analyst
Pyrene	BDL	μ	g/Kg			330	12/04/95	12/04/95	AT
Pyridine	BDL	μ	g/Kg			330	12/04/95	12/04/95	AT
Toxaphene	BDL	μ	g/Kg			3300	12/04/95	12/04/95	AT
1,2,4-Trichlorobenzene =	BDL	μ	g/Kg			330	12/04/95	12/04/95	AT
2,4,5-Trichlorophenol	BDL	μ	g/Kg			330	12/04/95	12/04/95	AT
2,4,6-Trichlorophenol	BDL	μ	g/Kg			330	12/04/95	12/04/95	AT
Surrogates									
Nitrobenzene-d5	49	Min: 23		Max:	120				
2-Fluorobiphenyl	72	Min: 30		Max:	115				
Terphenyl-d14	67	Min: 18		Max:	137				
Phenol-d5	55	Min: 24		Max:	113				
2-Fluorophenol	65	Min: 25		Max:	121				
2,4,6-Tribromophenol	68	Min: 19		Max:	122				
Volatile Organics				8240\8260		_			
Acetone	BDL	μ	g/Kg			250	11/30/95	11/30/95	AT
Benzene	BDL	μ	g/Kg			10	11/30/95	11/30/95	AT
Bromodichloromethane	BDL	μ	g/Kg			10	11/30/95	11/30/95	AT
Bromoform	BDL	μ	g/Kg			10	11/30/95	11/30/95	AT
Bromomethane	BDL	μ	g/Kg			10	11/30/95	11/30/95	AT
2-Butanone	BDL	μ	g/Kg			250	11/30/95	11/30/95	AT
Carbon disulfide	BDL	μ	g/Kg			10	11/30/95	11/30/95	AT
Carbon tetrachloride	BDL	μ	g/Kg			10	11/30/95	11/30/95	AT
Chlorobenzene	BDL	μ	g/Kg			10	11/30/95	11/30/95	AT
Chlorodibromomethane	BDL	μ	g/Kg			10	11/30/95	11/30/95	AT
Chloroethane	BDL	μ	g/Kg			10	11/30/95	11/30/95	AT .
Chloroform	BDL	μ	g/Kg			10	11/30/95	11/30/95	AT
Chloromethane	BDL		g/Kg			10	11/30/95	11/30/95	AT
Dibromomethane	BDL		g/Kg			10	11/30/95	11/30/95	AT
1,4-Dichloro-2-butene	BDL		g/Kg			10	11/30/95	11/30/95	AT
Dichlorodifluoromethane	BDL		g/Kg			10	11/30/95	11/30/95	AT
1,1-Dichloroethane	8DL		g/Kg			10	11/30/95	11/30/95	AT
1,2-Dichloroethane	BDL		g/Kg			10	11/30/95	11/30/95	AT
1,1-Dichloroethene	BDL		g/Kg			10	11/30/95	11/30/95	AT
trans-1,2-Dichloroethene	BDL		g/Kg			10	11/30/95	11/30/95	AT
1,2-Dichloropropane	BDL	-	g/Kg			10	11/30/95	11/30/95	AT

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Attn: HARRY HARSH

SOUTHERN WASTE SERVICES, INC. 1619 MOYLAN ROAD PANAMA CITY BEACH, FL. 32407 (800)852-8878 Page 10 6 Dec 1995 Report J5-11-080-02 LAB ID. 82223/E82101

Farameter	Result	Units	Method De	et. Limit	Extracted	Analyzed	Analyst
cis-1,3-Dichloropropene	BDL	μg/Kg		10	11/30/95	11/30/95	AT
trans-1,3-Dichloropropene	BDL	μg/Kg		10	11/30/95	11/30/95	7A
Ethanol	BDL	μg/Kg		1000	11/30/95	11/30/95	AT
Ethylbenzene	BOL	#Q/Kg		10	11/30/95	11/30/95	AT
Ethyl methacrylate	BOL	μg/Kg		10	11/30/95	11/30/95	AT
2-Rexanone	BDL	<b>иа/К</b> ф		100	11/30/95	11/30/95	AT
Lodomethane	BDL	μg/Kg		10	11/30/95	11/30/95	AT
Methylene chloride	BDL	μg/Kg		10	11/30/95	11/30/95	AT
4-Methyl-2-pentanone	BDL	μg/Kg		100	11/30/95	11/30/95	AT
Styrene	BDL	иg/Kg		10	11/30/95	11/30/95	AT
1,1,2,2-Tetrachloroethane	<b>S</b> DL	μg/Kg		10	11/30/95	11/30/95	ΑT
Tetrachloroethene	BDL	μg/Kg		10	11/30/95	11/30/95	AT
Toluene	BDL	μg/Kg		10	11/30/95	11/30/95	AT
1,1,1-Trichloroethane	BDL	μg/Kg		10	11/30/95	11/30/95	AT
1,1,2-Trichtoroethane	BDL	μ <b>g/</b> Kg		10.	11/30/95	11/30/95	AT
Trichloroethene	SDL	μg/Kg		10	11/30/95	11/30/95	AT
Trichlorofluoromethane	BDL	μg/Kg		10	11/30/95	11/30/95	AT
1,2,3-Trichloropropane	BOL	#8/KB		10	11/30/95	11/30/95	AT
Vinyl chloride	BDL	μg/Kg		10	11/30/95	11/30/95	AT
Xylenes (Total)	BDL	μg/Kg		10	11/30/95	11/30/95	AT
Surrogates							
Taluene-d8	95	Min: 81	Max: 117				
4-Bromofluorobenzene	89	Min: 74	Max: 121				
1,2-Dichloroethane-d4	95	Min: 70	Max: 121				

The D. Johnson, Valo Director

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Attn: HARRY MARSH

SOUTHERN WASTE SERVICES, INC. 1619 MOYLAN ROAD PANAMA CITY BEACH, FL. 32407 (800)852-8878 Page 11 6 Dec 1995

Report J5-11-080-03 LAB ID. 82223/E82101

Sample Description:

NCSC/JLS CONTRACTOR PANAMA CITY BEACH, FL.

COMPOSITE SAMPLE EXCAVATED SOILS

P.O. #: HM 95-2570

SAMPLE ID.: SAMPLE 003

COLLECTED: 11/28/95 RECEIVED: 11/28/95 COLLECTED BY: J. COOEY

Parameter	Result	Units	Method	Det. Limit	Extracted	Analyzed Analyst
ICAP Metals, TCLP			1311\6010			
Silver	BDL	mg/L		0.010	12/01/95	12/05/95 JC
Barium	0.078	mg/L		0.003	12/01/95	12/05/95 JC
Cadmium	BDL	mg/L		0.005	12/01/95	12/05/95 JC
Chromium	BDL	mg/L		0.005	12/01/95	12/05/95 JC
Lead	0.280	mg/L		0.100	12/01/95	12/05/95 JC
Arsenic	BDL	mg/L		0.075	12/01/95	12/05/95 JC
Selenium	BDL	mg/L		0.100	12/01/95	12/05/95 JC
Mercury, TCLP	<0.01	mg/L	1311\7471	0.01	12/01/95	12/01/95 JC
Purgeable Organics TCLP			1311\8260			
Benzene, TCLP	BDL	mg/L		0.20	11/30/95	11/30/95 AT
Carbon tetrachloride, TCLP	BDL	mg/L		0.20	11/30/95	11/30/95 AT
Chlorobenzene, TCLP	BDL	mg/L		0.20	11/30/95	11/30/95 AT
Chloroform, TCLP	BDL	mg/L		0.20	11/30/95	11/30/95 AT
1,4-Dichlorobenzene, TCLP	BDL	mg/L		0.20	11/30/95	11/30/95 AT
1,2-Dichloroethane, TCLP	BDL	mg/L		0.20	11/30/95	11/30/95 AT
1,1-Dichloroethene, TCLP	BDL	mg/L		0.20	11/30/95	11/30/95 AT
Methyl Ethyl Ketone, TCLP	BDL	mg/L		2.0	11/30/95	11/30/95 AT
Tetrachloroethene, TCLP	BDL	mg/L		0.20	11/30/95	11/30/95 AT
Trichloroethene, TCLP	BDL	mg/L		0.20	11/30/95	11/30/95 AT
Vinyl chloride, TCLP	BDL	mg/L		0.20	11/30/95	11/30/95 AT .
Surrogates						
4-Bromofluorobenzene	89	Min: 86	Max:	115		
Toluene-d8	97	Min: 88	Max:	110		
1,2-Dichloroethane-d4	101	Min: 76	Max:	114		•

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Attn: HARRY MARSH

SOUTHERN WASTE SERVICES, INC. 1619 MOYLAN ROAD PANAMA CITY BEACH, FL. 32407 (800)852-8878 Page 12 6 Dec 1995 Report J5-11-080-03 LAB ID. 82223/E82101

Greg D. Johnson, Lab Director

GLENT NAME: HE STY MOSSA	057 N. ELLIS ROAD, SUIT 908A BRECKENRIDGE P	TE 17, JACKSON' PARKWAY, TAMPA	/li 322 , FL 35510-42	254-2249 • (1 237 • (813) 62	904) 786- 26-0101	8340			30	J51	108	7		
11011	1	PROJECT NAME: A	Saval Ro	sporch L	35		PRESE	FINATIVE	5/	T	5/	1	77	$\overrightarrow{J}$
SWS Environmental  ADDRESS: 1609 Maylan Rd  Panama City Brac  PHONE:  FA	L, F1.	P.O. NUMBER / PRO	HID 95	5-75? City	Sito Brack	other	CONTA SIZE AND TYPE	MER 3	<i> -</i>	1		//		<del> </del>
CONTACT:  Marsh  TURN AROUND TIME OF RESULTS DUE BY  ENSTANDARD  D  RUSH  TO RUSH  T	X:	SAMPLED BY: JC   SPECIAL INSTRUCTI	lin C.				CRALLY ORDER D	8270	500	KA Metals			<i>t</i>	
SAMPLE	SAMPLE DESCRIPTION	- ** **	<del></del>	PLING	•	NO OF		+	$\bigcap$		+		LAE	8 USE
	ter separator Soil Sample		11/2 <b>8</b> /99	10/5	MATRIX	1	V	~ ~	~					
·						-								
											+			
* GWGroundwater SWSu	riace Water DW-Dri	Inking Water	WW-Waster	water	SO-Solid	d/Soil )		SLSlud			V_Ha	Indous W	hate	
FIELD PARAMETERS / COMMENTS:	217-011	ming Heler	7713310		INSFERS A	ELINOUISI				EPTED BY			130/95	A-A TIME 09:35
CONTUNERSISEALS INTACT	ONICE/4°C	SHIPPED	VIA	4	V			+		<del></del>	<u></u>	+	-	

P.002

TX/RX NO.1447

09:01

11/29/95

### Geos Laboratories Inc.

CORPORATE OFFICES 1057 N. ELLIS ROAD, SUITE 17 JACKSONVILLE, FL 32254-2249

(904) 786-8340 (800) 770-4367 (GEOS) FAX: (904) 786-7489

COLOGICAL, ENVIRONMENTAL AND OCEANOGRAPHIC SCIENCES

ANALYTICAL LABORATORY 1627 EAST 8th STREET JACKSONVILLE, FL 32206

(904) 354-6755 FAX: (904) 354-3799

S0U02

Attn: HARRY MARSH

SOUTHERN WASTE SERVICES, INC. 1619 MOYLAN ROAD PANAMA CITY BEACH, FL. 32407 Page 1 6 Dec 1995

Report J5-11-087-01 LAB ID. 82223/E82101

Sample Description:

OIL WATER SEPARATOR ADJACENT SOIL SAMPLES

LOCATION: PANAMA CITY BEACH/ NAVAL RESEARCH LABS

P.O. # HM 95-2570

SAMPLE ID.: #4

COLLECTED: 11/28/95 10:15
RECEIVED: 11/30/95
COLLECTED BY: J. COOEY

Parameter	Result	Units	Method	Det. Limit	Extracted	Analyzed	Analyst
Hydrocarbons, Total IR	12	mg/kg	9073	0.2	12/04/95	12/05/95	AM
ICAP Metals, Total			3050/6010				
Arsenic	BDL	mg/kg		7.5	12/05/95	12/06/95	JC
Barium	26	mg/kg		0.3	12/05/95	12/06/95	JC
Cadmium	1.0	mg/kg		0.5	12/05/95	12/06/95	JC
Chromium	9.2	mg/kg		0.5	12/05/95	12/06/95	JC .
Lead	976	mg/kg		10.0	12/05/95	12/06/95	JC
Selenium	BDL	mg/kg		10.0	12/05/95	12/06/95	1C
Silver	3.8	mg/kg		1.0	12/05/95	12/06/95	1C
Mercury, Total	0.104	mg/kg	7471	0.075	12/01/95	12/01/95	1C
Semi-Volatile Organics			3550\8270				
Acenaphthene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Acenaphthylene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Aldrin	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Aniline	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Anthracene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Benzidine	BDL	μg/Kg		660	12/04/95	12/04/95	AT
Benzoic acid	BDL	μg/Kg		1700	12/04/95	12/04/95	AT
Benzo (a) anthracene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Benzo (b) fluoranthene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Benzo (k) fluoranthene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Benzo (g,h,i) perylene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Benzo (a) pyrene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Benzyl alcohol	BDL	μg/Kg		330	12/04/95	12/04/95	AT
a-BHC	BDL	μg/Kg		330	12/04/95	12/04/95	AT
b-BHC	BDL	μg/Kg		330	12/04/95	12/04/95	AT
d-BHC	BDL	μg/Kg		330	12/04/95	12/04/95	AT
g-BHC (Lindane)	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Bis (2-chloroethoxy) methane	BOL	μg/ <b>Kg</b>		330	12/04/95	12/04/95	AT
Bis (2-chloroethyl) ether	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Bis (2-chloroisopropyl) ether	BDL	μg/Kg		330	12/04/95	12/04/95	AT

S0U02

Attn: HARRY MARSH

SOUTHERN WASTE SERVICES, INC. 1619 MOYLAN ROAD PANAMA CITY BEACH, FL. 32407 Page 2 6 Dec 1995 Report J5-11-087-01 LAB ID. 82223/E82101

arameter	Result	Units	Method	Det. Limit	Extracted	Analyzed	Analyst
Bis (2-ethylhexyl) phthalate	BDL	μg/Kg		330	12/04/95	12/04/95	AT
4-Bromophenyl phenyl ether	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Butylbenzylphthalate	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Chlordane	BDL	μg/Kg		1700	12/04/95	12/04/95	AT
4-Chloroaniline	BDL	μg/Kg		330	12/04/95	12/04/95	AT
4-Chloro-3-methylphenol	BDL	μg/Kg		330	12/04/95	12/04/95	AT
2-Chloronaphthalene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
2-Chlorophenol	BDL	μg/Kg		330	12/04/95	12/04/95	AT
4-Chlorophenyl phenyl ether	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Chrysene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
4,4'-DDD	BDL	μg/Kg		330	12/04/95	12/04/95	AT
4,41-DDE	BDL	μg/Kg		330	12/04/95	12/04/95	AT
4,4'-DDT	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Dibenzo (a,h) anthracene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Dibenzofuran	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Di-n-butylphthalate	BOL	μg/Kg		330	12/04/95	12/04/95	AT
1,2-Dichlorobenzene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
1,3-Dichlorobenzene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
1,4-Dichlorobenzene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
3,3'-Dichlorobenzidine	BDL	μg/Kg		660	12/04/95	12/04/95	AT
2,4-Dichlorophenol	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Dieldrin	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Diethylphthalate	BDL	μg/Kg		330	12/04/95	12/04/95	AT
2,4-Dimethylphenol	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Dimethylphthalate	BDL	μg/Kg		330	12/04/95	12/04/95	AT .
2,4-Dinitrophenol	BDL	μg/Kg		1700	12/04/95	12/04/95	AT
2,4-Dinitrotoluene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
2,6-Dinitrotoluene	BOL	μg/Kg		330	12/04/95	12/04/95	AT
Di-n-octylphthalate	BDL	μg/Kg		330	12/04/95	12/04/95	AT
1,2-Diphenylhydrazine	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Endosulfan I	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Endosulfan II	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Endosulfan sulfate	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Endrin	. BDL	μg/Kg		330	12/04/95	12/04/95	AT
Endrin aldehyde	BDL	μg/Kg		330	12/04/95	12/04/95	AT

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Attn: HARRY MARSH

SOUTHERN WASTE SERVICES, INC. 1619 MOYLAN ROAD PANAMA CITY BEACH, FL. 32407 Page 3 6 Dec 1995 Report J5-11-087-01 LAB ID. 82223/E82101

arameter	Result	Units	Method	Det. Limit	Extracted	Analyzed	Analyst
Fluoranthene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Fluorene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Heptachlor	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Heptachlor epoxide _	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Hexach Lorobenzene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Hexachlorobutadiene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Hexachlorocyclopentadiene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Hexachloroethane	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Indeno (1,2,3-c,d) pyrene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Isophorone	BDL	μg/Kg		330	12/04/95	12/04/95	AT
2-Methyl-4,6-dinitrophenol	BDL	μg/Kg		1700	12/04/95	12/04/95	AT
1-Methylnaphthalene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
2-Methylnaphthalene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
2-Methylphenol	BDL	μg/Kg		330	12/04/95	12/04/95	AT
4-Methylphenol	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Naphthalene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
2-Nitroaniline	BDL	μg/Kg		1700	12/04/95	12/04/95	AT
3-Nitroaniline	BDL	μg/Kg		1700	12/04/95	12/04/95	AT
4-Nitroaniline	BDL	μg/Kg		1700	12/04/95	12/04/95	AT
Nitrobenzene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
2-Nitrophenol	BDL	μg/Kg		330	12/04/95	12/04/95	AT
4-Nitrophenol	BDL	μg/Kg		1700	12/04/95	12/04/95	AT
N-Nitrosodimethylamine	BDL	μg/Kg		330	12/04/95	12/04/95	AT
N-Nitrosodiphenylamine	BDL	μg/Kg		330	12/04/95	12/04/95	AT
N-Nitrosodi-n-propylamine	BDL	μg/Kg		330	12/04/95	12/04/95	AT .
PCB-1016	BDL	μg/Kg		1700	12/04/95	12/04/95	AT
PCB-1221	BDL	μg/Kg		1700	12/04/95	12/04/95	AT
PCB-1232	BDL	μg/Kg		1700	12/04/95	12/04/95	AT
PCB-1242	BOL	· μg/Kg		1700	12/04/95	12/04/95	AT
PCB-1248	BDL	μg/Kg		1700	12/04/95	12/04/95	AT
PCB-1254	BDL	μg/Kg		1700	12/04/95	12/04/95	AT
PCB-1260	BDL	μg/Kg		1700	12/04/95	12/04/95	AT
Pentachlorophenol	BDL	. μg/Kg		1700	12/04/95	12/04/95	AT
Phenanthrene	BDL	μg/Kg		330	12/04/95	12/04/95	AT
Phenol	BDL	μg/Kg		330	12/04/95	12/04/95	AT

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Attn: HARRY MARSH

SOUTHERN WASTE SERVICES, INC. 1619 MOYLAN ROAD PANAMA CITY BEACH, FL. 32407 Page 4 6 Dec 1995 Report J5-11-087-01 LAB ID. 82223/E82101

Parameter	Result			Units	Method	Det	. Limit	Extracted	Analyzed	Analyst
Pyrene	BDL			μg/Kg			330	12/04/95	12/04/95	AT
Pyridine	BDL			μg/Kg			330	12/04/95	12/04/95	AT
Toxaphene	BDL			μg/Kg			3300	12/04/95	12/04/95	AT
1,2,4-Trichlorobenzene	BDL			μg/Kg			330	12/04/95	12/04/95	AT
2,4,5-Trichlorophenol	BDL			μg/Kg			330	12/04/95	12/04/95	AT
2,4,6-Trichlorophenol	BDL			μg/Kg			330	12/04/95	12/04/95	AT
Surrogates										
Nitrobenzene-d5	62	Min:	23		Max:	120				
2-Fluorobiphenyl	74	Min:	30		Max:	115				
Terphenyl-d14	70	Min:	18		Max:	137				
Phenol-d5	56	Min:	24		Max:	113				
2-Fluorophenol	73	Min:	25		Max:	121				
2,4,6-Tribromophenol	73	Min:	19		Max:	122				
Volatile Organics					8240\8260					
Acetone	BDL			μg/Kg			250	11/30/95	11/30/95	AT
Benzene	BDL			μg/Kg			10	11/30/95	11/30/95	AT
Bromodichloromethane	BDL			μg/Kg			10	11/30/95	11/30/95	AT
Bromoform	BDL.			μg/Kg			10	11/30/95	11/30/95	AT
Bromomethane	BDL			μg/Kg			10	11/30/95	11/30/95	AT
2-Butanone	BDL			μg/Kg			250	11/30/95	11/30/95	AT
Carbon disulfide	BDL			μg/Kg			10	11/30/95	11/30/95	AT
Carbon tetrachloride	BDL			μg/Kg			10	11/30/95	11/30/95	AT
Chlorobenzene	BDL			μg/Kg			10	11/30/95	11/30/95	AT
Chlorodibromomethane	BDL			μg/Kg			10	11/30/95	11/30/95	AT
Chloroethane	BDL			μg/Kg			10	11/30/95	11/30/95	AT .
Chloroform	BDL			μg/Kg			10	11/30/95	11/30/95	AT
Chloromethane	BDL			μg/Kg			10	11/30/95	11/30/95	AT
Dibromomethane	BDL			μg/Kg			10	11/30/95	11/30/95	AT
1,4-Dichloro-2-butene	BDL			μg/Kg			10	11/30/95	11/30/95	AT
Dichlorodifluoromethane	BDL			μg/Kg			10	11/30/95	11/30/95	AT
1,1-Dichloroethane	BOL			μg/Kg			10	11/30/95	11/30/95	AT
1,2-Dichloroethane	BDL			μg/Kg			10	11/30/95	11/30/95	AT
1,1-Dichloroethene	BDL			μg/Kg			10	11/30/95	11/30/95	AT
trans-1,2-Dichloroethene	BDL			μg/Kg			10	11/30/95	11/30/95	AT
1,2-Dichloropropane	BDL			μg/Kg			10	11/30/95	11/30/95	AT

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Attn: HARRY MARSH

SOUTHERN WASTE SERVICES, INC. 1619 MOYLAN ROAD PANAMA CITY BEACH, FL. 32407 Page 5 6 Dec 1995 Report J5-11-087-01 LAB ID. 82223/E82101

Parameter	Result	Units	Method D	et. Limit	Extracted	Analyzed	Analyst
cis-1,3-Dichloropropene	BDL	μg/Kg		10	11/30/95	11/30/95	AT
trans-1,3-Dichloropropene	BDL	μg/Kg		10	11/30/95	11/30/95	AT
Ethanol	BDL	μg/Kg		1000	11/30/95	11/30/95	AT
Ethylbenzene	BDL	μg/Kg		10	11/30/95	11/30/95	AT
Ethyl methacrylate	BDL	μg/Kg	•	10	11/30/95	11/30/95	AT
2-Hexanone	BOL	μg/Kg		100	11/30/95	11/30/95	AT
Iodomethane	BDL	μg/Kg		10	11/30/95	11/30/95	AT
Methylene chloride	BDL	μg/Kg		10	11/30/95	11/30/95	AT T
4-Methyl-2-pentanone	BDL	μg/Kg		100	11/30/95	11/30/95	AT
Styrene	BDL	μg/Kg		10	11/30/95	11/30/95	AT
1,1,2,2-Tetrachloroethane	8DL	μg/Kg		10	11/30/95	11/30/95	AT
Tetrachloroethene	BDL	μg/Kg		10	11/30/95	11/30/95	AT
Toluene	BDL	μg/Kg		10	11/30/95	11/30/95	AT
1,1,1-Trichloroethane	BDL	μg/Kg		10	11/30/95	11/30/95	AT
1,1,2-Trichloroethane	BDL	μg/Kg		10	11/30/95	11/30/95	AT
Trichloroethene	BDL	μg/Kg		10	11/30/95	11/30/95	AT
Trichlorofluoromethane	BDL	μg/Kg		10	11/30/95	11/30/95	AT
1,2,3-Trichloropropane	BDL	μg/Kg		10	11/30/95	11/30/95	AT
Vinyl chloride	BDL	μg/Kg		10	11/30/95	11/30/95	AT
Xylenes (Total)	BDL	μg/Kg		10	11/30/95	11/30/95	AT
urrogates							
Toluene-d8	99	Min: 81	Max: 117				
4-Bromofluorobenzene	102	Min: 74	Max: 121				
1,2-Dichloroethane-d4	107	Min: 70	Max: 121				

dreg D. Johnson Jab Director

"DDY	RECO	)RD

1057 N ELLIS ROAD, SUITE 17, JACKSONVILLE, FL 32254-2249 • (904) 786-8340 5909A BRECKENRIDGE PARKWAY, TAMPA, FL 33610-4237 • (813) 626-0101

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Para City Brack Fl.	PROJECT LOCATION:	111.35 - Area #	2570 333 US)	T site		CONTE SIZE AND TYPE	Kur /		//	<u> </u>		//	
CONTACT Harry Marsh	SAMPLED BY:  Tulia  SPECIAL INSTRUCTION	: C. C		1		¢nmo <r>&gt;&gt; ≥</r>	T /						
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MW-1 Crundwater, UST area # 33	3	12-15-95	10:10	6w	3	~	1	1	-51			\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\	
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## Geos Laboratories Inc.

CORPORATE OFFICES
1057 N. ELLIS ROAD, SUITE 17
JACKSONVILLE, FL 32254-2249

(904) 786-8340 (800) 770-4367 (GEOS) FAX: (904) 786-7489

GEOLOGICAL, ENVIRONMENTAL AND OCEANOGRAPHIC SCIENCES

ANALYTICAL LABORATORY 1627 EAST 8th STREET JACKSONVILLE, FL 32206

(904) 354-6755 FAX: (904) 354-3799

S0U02

Attn: HARRY MARSH

SOUTHERN WASTE SERVICES, INC. 1619 MOYLAN ROAD PANAMA CITY BEACH, FL. 32407 Page 1 22 Dec 1995

Report J5-12-049-01 LAB ID. 82223/E82101

Sample Description: UST OIL/WATER SEPARATOR

AREA #333 UST SITE/USCG-NAS PANAMA CITY BEACH, FL.

GROUNDWATER

SAMPLE ID.: MW-1

COLLECTED: 12/15/95 10:10

RECEIVED: 12/18/95
COLLECTED BY: J. COOEY

Parameter	Result	Units	Method	Det	. Limit	Extracted	Analyzed	Analyst
Hydrocarbons, Total IR	1.79	mg/L	418.1		0.200	12/21/95	12/21/95	AM
Polynuclear Aromatics			625\8270					
Naphthalene	BDL	μg/L		,	10	12/20/95	12/21/95	AT
Acenaphthylene	BDL	μg/L			10	12/20/95	12/21/95	AT
1-Methylnaphthalene	49	μg/L			10	12/20/95	12/21/95	TA
2-Methylnaphthalene	BDL	μg/L			10	12/20/95	12/21/95	AT
Acenaphthene	BDL	μg/L			10	12/20/95	12/21/95	AT
Fluorene	BDL	μg/L			10	12/20/95	12/21/95	AT
Phenanthrene	BDL	μg/L			10	12/20/95	12/21/95	AT
Anthracene	BDL	μg/L			10	12/20/95	12/21/95	AT
Fluoranthene	BDL	μg/L	•		10	12/20/95	12/21/95	AT
Pyrene	BDL	μg/L			10	12/20/95	12/21/95	AT
Benzo(a)anthracene	BDL	μg/L			10	12/20/95	12/21/95	AT
Chrysene	· BDL	μg/L			10	12/20/95	12/21/95	AT
Benzo(b)fluoranthene	BDL	μg/L			10	12/20/95	12/21/95	AT
Benzo(k)fluoranthene	BDL	μg/L			10	12/20/95	12/21/95	AT
Benzo(a)pyrene	BDL	μg/L			10	12/20/95	12/21/95	ΑT
Indeno(1,2,3-c,d)pyrene	BDL	μg/L			. 10	12/20/95	12/21/95	ΑT
Dibenzo(a,h)anthracene	BDL	μg/L			10	12/20/95	12/21/95	AT
Benzo(g,h,i)perylene	BDL	μg/L			10	12/20/95	12/21/95	AT
Surrogates			• -					
Nitrobenzene-d5	74	Min: 35	Max	114				•
2-Fluorobiphenyl	80	Min: 43	Max	116				
4-Terphenyl-d14	50	Min: 33	Max	141				
Volatile Aromatics	•		602					
Methyl-tert-butyl ether	BDL	μg/L			5.0	12/18/95	12/18/95	OLS
Benzene	1.5	μg/L			1.0	12/18/95	12/18/95	OLS
Toluene	BDL	μg/L			1.0	12/18/95	12/18/95	OLS
Ethyl benzene	1.0	μg/L			1.0	12/18/95	12/18/95	OLS
Xylene, Total	5.2	μg/L			1.0	12/18/95	12/18/95	OLS
Chlorobenzene	BDL	μg/L			1.0	12/18/95	12/18/95	OLS

SOUGE

Attn: - .. BARSH

SOUTHER FASTE SERVICES, INC.

1619 -- 1 - 20 ACAD

PANAHA 3EACH, FL. 32407

Page 2 2Z Dec 1995 Report J5-12-049-01 LAB ID. 82223/E82101

Paramer::	Result	Units Het	hod Det.	Limit	Extracted	Analyzed	Anelyst
1,4-1 shippropensene	90L	μg/L		1.0	12/18/95	12/18/95	OL\$
1,3 4 m torobenzene	BOL	#9/L		1.0	12/18/95	12/18/95	OLS
1,2 Prorobenzene	BDL	μg/l		1.0	12/18/95	12/18/95	OLS
Surroj 11:3							
Brown with the	102 M	in: 70	Nex: 130				

Greg D. Johnson Lab Director

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# APPENDIX D SOIL BORINGS LOGS

LOG OF BORING

		1
SHEET	OF	1

LOC	ATIC	ON O	F BC	RIN	G:					•		PROJECT: CTO C	068	BORING NO. 5801				
													Site 333 TOTAL DEPTH: 4.5"					
												JOB NO.	7113	LOGGED BY: 6, ItUMS				
												PROJ.MGR:		EDITED BY:				
												DRILLING CONTRACTOR:	NA					
يح	و	5i ł	ٹ (	30	riv	19	Loc	atic	>in M	La	P	DRILL RIG TYPE:	NA					
						J						DRILLER'S NAME	NA					
				-								SAMPLING METHODS:	hand	auger				
												STARTED TIME:	1600	DATE: 6/10/96				
												COMPLETED TIME:	1639	DATE: 6/10/96				
												BORING DEPTH (fL)	4.51					
												CASING DEPTH (ft.)	NA					
												WATER DEPTH (ft.)						
SAMPLE DEPTH	子 SAMPLER TYPE	BLOWS/6-IN.	INCHES DRIVEN	INCHES RECOVERED	A کونی Moisture		UNFILTERED OVA (PPM)	(y) FILTERED OVA (PPM)	CORRECTED OVA (PPM)	DEPTH IN FEET	N X X	SAND, I+gy Fg~mg, s	ay, fredome silt,	ominantly atz. , loose, moist				
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LOG OF BORING

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LOC	CATIC	ON O	F BO	RIN	G:							PROJECT: CTO	०००८	BORING NO. 5BOZ			
												Site 333 TOTAL DEPTH: 4.5"					
ĺ												JOB NO.	7113	LOGGED BY: G. Helms			
												PROJ.MGR:		EDITED BY:			
												DRILLING CONTRACTOR:	NA				
5	ee.	50	i 1 1	B0	rir	la	しむこ	atio.	on 1	الانك	2	DRILL RIG TYPE:	NA				
						_						DRILLER'S NAME	NA				
												SAMPLING METHODS:	hand a	war			
												STARTED TIME:	1603	DATE: 6/10/98			
												COMPLETED TIME:	1637	DATE: 6/10/96			
												BORING DEPTH (ft.)	4.5'				
												CASING DEPTH (fL)	NA				
												WATER DEPTH (ft.)					
SAMPLE DEPTH	SAMPLER TYPE	BLOWS/6-IN.	INCHES DRIVEN	INCHES RECOVERED	MOISTURE	ODOR	UNFILTERED OVA (PPM)	FILTERED OVA (PPM)	CORRECTED OVA (PPM)	DEPTH IN FEET	USCS CODE						
<u>  '</u> _					<u> </u>	<u> </u>			<u> </u>	1 1 -	4						
2	нн	_		_	المرم		0	٥	0	2	-	SAND, It.	Ocad	ominanthiatz.			
$\vdash$					<u> </u>			-		1 1.7	54	) <u> </u>		loose some			
3										3 🗴	P''	pepple si	<del>,</del>	avavel moist.			
					NO		10	3	2		]			,			
4	144				74-					4	١.,	SI HY SAN	Sidk gra	4, organic rich,			
5									1		5M		race Debl				
F					-				<del></del>	5	1	graver, we	1 at 4,01.	Noted a layer			
6								1		6	1	below 4.0'	fa, high	sandy Sict, just			
											1	content.	3, 119	13011			
7					<u>L</u>					7	]						
8																	
13	ļ.,			<u> </u>		_	<u> </u>	ļ	<u> </u>	8							
9				ļ						-	1						
$\vdash$	$\vdash$				<u> </u>	-		<u> </u>	<u> </u>	9	1		<u> </u>				
-									1	-	1		- <del></del>				
NO.	- NO	ORC	SANI	C VA	L	I RS D	ETEC	L	L	1 L	1	SOIL/SEDIMENT DESCRIF	TION				

LOG OF BORING

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LOC	CATIC	о ис	FBC	RIN	G:							PROJECT: CTO C	CCS.	BORING NO. 5803
												Site ?	33	TOTAL DEPTH: 4.57
												JOB NO.	7113	LOGGED BY: 6, Items
												PROJ.MGR:	6.600de	EDITED 8Y:
		<i>~</i> :	١	D			٠ ١ ٥	~ ~ <del>\</del>				DRILLING CONTRACTOR:	NA	
) t	وو	21	He	10	OVI	irl	9 4	car	ion	Ma	P	DRILL RIG TYPE:	NA	
												DRILLER'S NAME	NA	
												SAMPLING METHODS:	hand a	uger
												STARTED TIME:	0803	DATE: 6/1/196
COMPLETE													0826	DATE: 6/11/96
												BORING DEPTH (ft.)	4.51	
												CASING DEPTH (ft.)	NA	
												WATER DEPTH (fL)		
	SAMPLER TYPE	BLOWS/6-IN.	INCHES DRIVEN	INCHES RECOVERED	MOISTURE	) ODOR	UNFILTERED OVA (PPM)	Q) FILTERED OVA (PPM)	り corrected ova (PPM)	T DEPTH IN FEET	USCS CODE	SAND, It go	ray, banc	ted widk gray Sand is predem
3										3 🔀	۷,۸	inantly 9.	by poor	graded,
4	HA	-	-	_	NET	_	29	10	19	_ _		moist fg	-mg.	
<u> </u>					7					<b>4</b> <del>X</del>	١,	SAND; dko	tray, to	ilt, poexietym
5										5 <u>X</u>	24	graded, w		
G										6		3		
7										7			<del></del>	
8										-				
├										8				
9										9				
						Н							<del></del>	
			-								1			
ND:	= NO	ORC	ANI	C VA	POR	25.0	ETECT	FD		·	•	SOIL/SEDIMENT DESCRIP	TION	

LOG OF BORING

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															<u> </u>			
LOC	CATIC	о ис	F BC	RIN	G:								PROJECT: CTO 0008 BORING NO. PCY -333-MWC					
ļ													Site	333	TOTAL DEPTH: 12.2			
													JOB NO.	7113	LOGGED BY: G. Helms			
													PROJ.MGR:	G. Goode	EDITED BY:			
												i	DRILLING CONTRACTOR: NA					
5	ee	5	·le	Bo	viv	rα	Loc	ah	oni	U,	ap		DRILL RIG TYPE: NA					
	See Site Boring Location Map												DRILLER'S NAME NA					
													SAMPLING METHODS:	hand o	walr			
													STARTED TIME:	0830	DATE: 611196			
COMPLETED TIME: 0908 DATE: 611190																		
													BORING DEPTH (ft.)	12.21				
													CASING DEPTH (ft.)	NA				
	_												WATER DEPTH (ft.)					
										_								
							Σ		<b>⊋</b>									
	}			ED			d.	Ž	9									
I	ш		z	VER			≸	ا ق	×	_		ı						
EPT	ΙΥΡ	ż	SIVE.	00	_		8	8	g	E	u	ų						
ED	ER	F9/9	S DF	S RE	JRE		ER	9	CT	Z	5	5						
SAMPLE DEPTH	SAMPLER TYPE	BLOWS/6-1N.	INCHES DRIVEN	INCHES RECOVERED	MOISTURE	ODOR	UNFILTERED OVA (PPM)	FILTERED OVA (PPM)	CORRECTED OVA (PPM)	DEPTH IN FEET	3000 SOSI	ŝ						
SAI	SA	ВГ	NC.	INC	ΨO	8	3	ᇤ	00	OE	=	3			·····			
,							•				_	1						
1										1								
2	HA			_	CUBY	_	2	0	2									
	7 7				6	$\vdash$				2	$\exists$	1	SAND, It. to	n, fg-mg	, predominantly			
3										_	XS	P	gtz. poor	y graded	predominantly			
_	-					-		-		3	4		med grai	near, trac	e silt, moist.			
4	HA	_			¥	-	2	0	2	4	-		SAND, It.a.	mus Vicus	fine to med.			
											$\nabla$	ı			hyatz. osorhy			
5	HA	_	_	_	NE	_	100	40	60	5	$\overline{\mathbf{x}}$	ŀ	aradea (50		care very fine			
											X		Some silt	Wet-41,5	,			
ķ										в	$\boxtimes$							
7				_	1		100	dia.	60				* 5,0' note	d sand an	d silt mixed			
1	HΛ	_			WE		100	40	~	7			digray	o blk, di	esel like coor			
8											$\overline{X}$		WIBILYSU	een				
, J										8	X		<u> </u>		1			
9											$\dashv$		* 1.0' - SAN	s as above	. W/ Strong dresel			
-										9	$\dashv$		odor.	1, 1, 0, 0, 0, 0	me material			
											$\dashv$		4 8.0 - 5H		The mount			
1	7											•	• • • • • • • • • • • • • • • • • • •	\/ ·				

ND = NO ORGANIC VAPORS DETECTED

NS = NO CARBON FILTERED SAMPLE READ

SOIL/SEDIMENT DESCRIPTION

\* PCY-333-MWOH was extended by hand, by drillers to 12.2' bls.

LOG OF BORING

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LOC	ATIC	O NC	F BC	RIN	G:							PROJECT:	CTO	०००४	BORING NO. 5805
													site	333	TOTAL DEPTH: 4.21
												JOB NO.		7113	LOGGED BY: G. Helms
												PROJ.MGR:		6.60de	
												DRILLING CONTI	RACTOR:	NA	
												DRILL RIG TYP	E:	NA	
												DRILLER'S NA	ME	NIA	
												SAMPLING ME	THOOS:	hand	wer
												STARTED TIM	E:	0955	DATE: 611196
												COMPLETED	TIME:	1031	DATE: 6/11/96
												BORING DEPT	H (fL)	4.21	
												CASING DEPT	H (ft.)	NA	
	,					,						WATER DEPTI	H (ft.)		
DOLD OF CONTROPTH	ま ま SAMPLER TYPE		I INCHES DRIVEN		A A A A MOISTURE	ODOR	D C UNFILTERED OVA (PPM)	D O FILTERED OVA (PPM)	O O CORRECTED OVA (PPM)	1 2 3 4 5 6 7 8 9		SAND GIZ: trace Somo gray May May Cabur	predo 5; 1+, ova	minantly well are anic mo color.  ith SANI poorly arganic	Jeriai, ak Didk gray, fg- graded,
Ė			-						<del> </del>	▎▝├	$\dashv$				
1					L										
				<u> </u>			ETEC					COU /SEDIMEN			

NS = NO CARBON FILTERED SAMPLE READ

SOIL/SEDIMENT DESCRIPTION

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LOG OF BORING

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LOC	ATIC	о ис	F BC	RIN	G:				-			PROJECT: CTO	0008	BORING NO. 5806
												site "	333	TOTAL DEPTH: 4.51
												JOB NO.	7113	LOGGED BY: G. Helms
1												PROJ.MGR:	<del></del>	EDITED BY:
[												DRILLING CONTRACTOR:	NA	
3	See soil Boring Location Map DRILL RIG TYPE: NA													
DRILLER'S NAME NT+														
İ	SAMPLING METHODS: hand auger													
												STARTED TIME:	1115 HAVM	DATE: 6/11/96
												COMPLETED TIME:	1149	DATE: 6/1/96
												BORING DEPTH (ft.)	4.51	
												CASING DEPTH (ft.)	NA	
												WATER DEPTH (ft.)		
							₩.		æ.		1	Į		
				RED			(P	₽	<u>4</u>	j	}	1		
Ξ	jĘ.		N.	)VE			ð	A (P	ð	<u></u>				. [
EP	ΤΥ	ż	RIV	ECC	l in		G.	8	ED	E	삙			
<u> </u>	LER	S/6.	S D	SR	2.		TER	GE .	ECT	Z ∓	Ö	1		1
SAMPLE DEPTH	SAMPLER TYPE	BLOWS/6-IN.	NCHES DRIVEN	INCHES RECOVERED	MOISTURE	ODOR	UNFILTERED OVA (PPM)	FILTERED OVA (PPM)	CORRECTED OVA (PPM)	DEPTH IN FEET	USCS CODE			
\s	's	10	Z	Z	Σ	ō	5	區	<u> </u>	٦	j			
1										⊢, ا	}			
			_		-			-		'}-	1			
2	HA	_	_	-	moth	-	0	0	0	2	1	SAND, dk gr	50 p 60 = 0	22 Oxedonional
					10°					×	1	9+7 50 M	silt. som	e organic
3	HA	-	-	-	WED.	-	4	4	4	3 🗙		material	maist.	9,90,
. :										区	1	17.15.7		
4										4 🗵	]	SAND It.	tan fa-r	ng, (predominantly
5												med avair	1, ' 4	ominantly gtz.
										5	1	trace silt,	trace or	ganicmourial,
6		,								<u> </u>	↓	moist fro		321 then wet
10					_	_		<u> </u>		6	1			rom 2,8-3,21
7											-	Continued	handai	sering to 4.5%
Ľ				<u> </u>		-		<del> </del>		7 _	-	Sand was	s Satura	diesel-like
8								1		_	┨	and had	a Strong	alesel -like
		-	-	-	-			<del>                                     </del>		8	1	odor.	<del> </del>	
9				1	[					<sub>e</sub>	1			
<del>-</del>			-	-	-	<del>                                     </del>		-		" -	1			<del></del>
			}	j						<u> </u>	1		· <u> </u>	
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ND = NO ORGANIC VAPORS DETECTED NS = NO CARBON FILTERED SAMPLE READ SOIL/SEDIMENT DESCRIPTION

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LOG OF BORING

LOC	OCATION OF BORING:											PROJECT: CTO CO	08	1004 BORING NO. 5B07
												5He 333	_	TOTAL DEPTH: 5
												JOB NO.	7//3	LOGGED BY: 6. 600de
												PROJ.MGR:	6.6.0de	EDITED BY: 6. 6000
		,										DRILLING CONTRACTOR:	NA	
/	•	ر. رونون			, .			_				DRILL RIG TYPE:	NA	
5	pe.	iΒ	oci.	ng	ho	ca.	tion	. 4	·P			DRILLER'S NAME	NA	
				J					•			SAMPLING METHODS:	Hond Ave	21
												STARTED TIME:	12:00	DATE: 6/11/96
												COMPLETED TIME:	12:15	DATE: 6/11/96
												BORING DEPTH (ft.)	51	
												CASING DEPTH (ft.)	NA	
												WATER DEPTH (ft.)	_	
Z SAMPLE DEPTH	SAMPLER TYPE	l BLOWS/6-IN.	INCHES DRIVEN	I WOHES RECOVERED	Decrease A	9CG0) 1	ONFILTERED OVA (PPM)	FILTERED OVA (PPM)	CORRECTED OVA (PPM)	DEPTH IN FEET	NSCS CODE	Sond, light gray trace of silt,	fice to me	dive grained quartz,
3										3 X	[	1005e, dry		
4	¥А	-	_	-	wet	-	مه	NS	ND	4		Sond light brow	in Fine t	medium grained,
5			!							5 X	SŒ	Some Silt of 10	plasticit	we tat 4.5"
			j		_							(oily sheen to	11.1	[ o( )
6					·					8	}	/	<del></del>	
										7	1			
											]			
										8	-			
										9	1			
											]			
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Permit #T9601496 Two4 SHELL OF

LOG OF BORING

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SHEET	OF	1

LO	CATIC	о ис	F BC	RIN	 G:								PROJECT: CTO	0005	BORING NO. 5808
													site	333	TOTAL DEPTH: 5,0/
													JOB NO.	7113	LOGGED BY: 6. Helms
													PROJ.MGR:	6.600de	EDITED BY:
													DRILLING CONTRACTOR:	NA	
													DRILL RIG TYPE:	NIA	
15	See Soil Boring Location Map												DRILLER'S NAME	NA	
l													SAMPLING METHODS:	hand au	ich
ļ													STARTED TIME:	1200	DATE: 6/11/96
													COMPLETED TIME:	1232	DATE: 6/1/96
													BORING DEPTH (fL)	5,0'	
													CASING DEPTH (ft.)	NA	
													WATER DEPTH (ft.)		
SAMPLE DEPTH	SAMPLER TYPE	BLOWS/6-IN.	INCHES DRIVEN		MOISTURE		UNFILTERED OVA (PPM)	O FILTERED OVA (PPM)	O CORRECTED OVA (PPM)	DEPTH IN FEET		USCS CODE	SAND, It. to		nches of concrete g (predominantly jtz., prorky
4	/+A		_	-	ROF VIE		0	0	0	4	$\frac{1}{}$			an, fg-mg	Coredominanthy
5										5	X.	R	med grains antly 9tz	of, trace :	silt, predomin-
Ы										в	$\exists$		to~4.0' 11	en wet	Sample taken
1											4		<del></del>	···-	
<u> </u>									<u></u>	7 -	4		* 5.0'-hi	+ saxar	15, CT; Fg,
as I										8			fuel like	city i SA	T, sily sheen,
a											$\perp$			<u> </u>	
-1									<u> </u>	9	4			<del></del>	
										-	$\dashv$				
	ليبا			2 1 / 4			ETEC		l	ı L			SOIL/SEDIMENT DESCRIP	TION	

LOG OF BORING

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												<i>ca</i> 0
LOC	LOCATION OF BORING:											PROJECT: CTO 0008 BORING NO. 333-TW06
	reaction of parities											Site 333 TOTAL DEPTH: 6.0
												JOB NO. 7113 LOGGED BY: G. Helms
												PROJ.MGR: G. GOODE EDITED BY:
												DRILLING CONTRACTOR: NA
_	see soil boring Location Map											DRILL RIG TYPE: NA
J	حر	- =	×Οι	10	۱ ۷ ت	V V	5 4	عرم	1701	i i ota	Т,	DRILLER'S NAME NA
												sampling methods: hand auger
												STARTED TIME: 1400 DATE: 611196
												COMPLETED TIME: 1426 DATE: 611196
												BORING DEPTH (ft.)
												CASING DEPTH (R.)
L												WATER DEPTH (ft.)
SAMPLE DEPTH	SAMPLER TYPE	BLOWS/6-IN.	NCHES DRIVEN	NCHES RECOVERED	MOISTURE	ODOR	UNFILTERED OVA (PPM)	FILTERED OVA (PPM)	CORRECTED OVA (PPM)	DEPTH IN FEET	USCS CODE	
	٠,		=	=		Ü					-	Cored through 6 inches of Concreto
1					_					1 _		before hand augering.
~	منــن				not y	_	-	_	٦.			,
2	İΤΑ				60		3	9	7_	2		SAND, it. gray, fg-mg gtz., some
3										3 X		silt, Wellanded, moist, some or- ganic material (dkgray) throughout
4	HA		_	_	WE	_	5	2	3	4		SAND, It gray, fg-mg giz, predom-
5					4					又		inanthemed arained poorly
7										5 X		sorted trace organic material
(c)												wet,
							<u>.</u>			6		
7					1					_	l	
						$\vdash \vdash$				7		* 4.5'-5.0' - 011 sheen in 50:15
5				}						⊢ا。 ا	1	Was fine amined hore bowersh
					<u> </u>	<b></b>			L	8	l	was fine grained here, brownish

ND = NO ORGANIC VAPORS DETECTED
NS = NO CARBON FILTERED SAMPLE READ

SOIL/SEDIMENT DESCRIPTION

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\* previously SB09

LOG OF BORING

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LO	CATIO	о ис	FBC	ORIN	G:							PROJECT: CTO-0008 BORING NO. 333-TWO				
												si te	333	TOTAL DEPTH: 6.07 40		
ľ												JOB NO.	7113	LOGGED BY: G, Itelms		
ļ												PROJ.MGR:	6,000di	EDITED BY:		
												DRILLING CONTRACTOR:	NA			
50	, ,	50	( ) 1	Ro-	ンハ	na	L00	cati	on	Ma	D	DRILL RIG TYPE:	NA			
			•	<i>-</i>	• •	5	)				7	ORILLER'S NAME	NA			
												SAMPLING METHODS:	hand	waer		
												STARTED TIME:	1620	DATE: 6/11/96		
												COMPLETED TIME:	1652	DATE: 6111196		
												BORING DEPTH (ft.)	5,01			
1												CASING DEPTH (ft.)	NA			
												WATER DEPTH (ft.)				
											$T^-$					
}	j						€	1	_ <b>∑</b>		1					
	ļ			ED			dd)	ξ	9							
I	ш		2	ER	}		. ₹	a a	≸	_						
EP T	<u>ح</u>	z	3	000	_			8	ä	133						
	ER	11-9/	P.O.	RE.	JRE	]	ERE	i ii	2	3	CODE					
SAMPLE DEPTH	SAMPLER TYPE	BLOWS/6-1N.	INCHES DRIVEN	INCHES RECOVERED	MOISTURE	8	UNFILTERED OVA (PPM)	FILTERED OVA (PPM)	CORRECTED OVA (PPM)	DEPTH IN FEET	SS					
SAI	SAI	BLC	2	2	₩ W	ODOR	Š	글	00	DEI	USCS					
										ΙL	]	Coved throw	an ~ bir	ches of		
								L		1	_	Concrete 0	rior to b	oring.		
٦	امنا				mor	1_		_			_			<u> </u>		
2	$H\Delta$	_	_		40°		0	0	0	2		SAND, It. tau	n, fg-mg	predominantly		
2										2	358	gtz., Wson	e organic	material		
3	_				├					3	4	Kakgran +	oblks, pre	dominantly		
4	HA		_	_	JET		0	0	0	-	-	med grain	ed, poorly	sorted, moist.		
<del>                                     </del>	-			-	3	$\vdash$				1 4		SAND, It go	cas to m	3 atz. some		
5					}					5 \	75	organic m	naturally	Well graded,		
<u> </u>			_		_	$\vdash$		<del> </del>		"	4			stom of sample		
4				Ì						6	1	Jecha 13 A	10W1 + C	Bur fer by tubres.		
<u> </u>					$\vdash$						1	5,01-Sar	d Cine	grained inced		
7										7	7	antimed.	It. aran	trace silt,		
						П				1	1	proanic o	maderial			
S										8	]	Well sorted		WIFID.		
<u></u>																
9									<u></u>	9	1					
1	] ,							,		L	_					
L	لبا	لـا				Ш				l L	╛					
MO	- 110	000	TAALI.	~ \ / A		200	ETECT					COLUCEDIMENT DESCRIPT	TION			

\* Permit # T9601494

NS = NO CARBON FILTERED SAMPLE READ

LOG OF BORING

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100	ATIC	N O	F BO	RING	G:							PROJECT: CTO	0008	BOBING NO :322 -11-2
	,,,,,												333	BORING NO. 333 TWO3
}												JOB NO.	7113	
												PROJ.MGR:	G. Goode	LOGGED BY: G. Helms
												<del></del>	NA NA	JEDITED BY:
5	90	So	NE	301	in	a	LOC	io:	on f	ron	?	DRILLING CONTRACTOR:	NA	
						-						DRILL RIG TYPE:	NA	
												DRILLER'S NAME	<del> </del>	
ļ												SAMPLING METHODS:	T	auger
												STARTED TIME:	1700	DATE: 6/1/96
												COMPLETED TIME:	1722	DATE: 6/11/96
												BORING DEPTH (ft.)	5,01	
												CASING DEPTH (R.)	NA	
		<del></del>				,		1				WATER DEPTH (fL)		
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ND = NO ORGANIC VAPORS DETECTED NS = NO CARBON FILTERED SAMPLE READ

\* Permit # T9 601495 \* SBII (previous boring)

SOIL/SEDIMENT DESCRIPTION

LOG OF BORING

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												5 te 3	33	TOTAL DEPTH: 5/11 615
												JOB NO.	7113	LOGGED BY: G. 1. LLIMS
												PROJ.MGR:	G. Goode	EDITED BY:
												DRILLING CONTRACTOR:	NA	
				_				c a +		Ma	$\cap$	DRILL RIG TYPE:	NA	
5	e°	50	<i>i</i> 1	Bo	γŶΥ	رم	LO	انص	ion	1000	Α-	DRILLER'S NAME	NA	
												SAMPLING METHODS:	handa	lacer
												STARTED TIME:	0955	DATE: 6/12/95
												COMPLETED TIME:	0925	DATE: 6/12/95
												BORING DEPTH (ft.)		
												CASING DEPTH (fL)	NA	
Ì												WATER DEPTH (ft.)		
S 2 9 1 5 1 5 AMPLE DEPTH	(大) (大) (大) (大) (大) (大) (大) (大) (大) (大)	BLOWS/6-IN.	INCHES DRIVEN	I INCHES RECOVERED	A MOISTURE,		() O UNFILTERED OVA (PPM)	U O FILTERED OVA (PPM)	O CORRECTED OVA (PPM)	1 2 3 4 5 6 7 8 1 DEPTH IN FEET	USCS CODE	prior to he DAND, It. predominar fg-mg; m	tan wy sor tan wy sor with 9 +7., oisi in, fg-me graded, r sample w	ne brown grains little silt, predominantly noist to a 3:8' wet
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NS = NO CARBON FILTERED SAMPLE READ

<sup>\*</sup> Permit # T9601497

<sup>\*</sup> TWOS was previously SB12.

LOG OF BORING

														PC4 - 333 - MW
LOC	CATIC	O NC	FBC	RIN	G:							PROJECT: CTO	0003	BORING NO.
												site	333	TOTAL DEPTH: 12.0
												JOB NO.	7/13	LOGGED BY: G. Helms
												PROJ.MGR:	G. Goode	EDITED BY:
5	عد											DRILLING CONTRACTOR:	Groundu	later Protection
51	re	50	oil	b	ori	n	z Lo	cat	ior	Mo	4	DRILL RIG TYPE:	D120C	(Diedrick)
											•	DRILLER'S NAME	charles	Bucher
												SAMPLING METHODS:	hand au	ger / Spij + Speci
												STARTED TIME:	1713	DATE: 6/12/96
												COMPLETED TIME:		DATE:
												BORING DEPTH (ft.)		
												CASING DEPTH (ft.)	NA	
		<u> </u>			<b>,</b> .	,—.				,		WATER DEPTH (ft.)		
SAMPLE DEPTH	SAMPLER TYPE	BLOWS/6-IN.	INCHES DRIVEN	INCHES RECOVERED	MOISTURE	ODOR	O UNFILTERED OVA (PPM)	O FILTERED OVA (PPM)	CORRECTED OVA (PPM)	DEPTH IN FEET		SAND, brow antly mg,	on, fa-r	ng predomin- graded; qtz.
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SOIL/SEDIMENT DESCRIPTION

NS = NO CARBON FILTERED SAMPLE READ

ND = NO ORGANIC VAPORS DETECTED

\* previously SBIO.

ü = DEPTH TYPE (b) BLOWS DRIVEN - CEC'VD' MET MOISTURE ODOR 0 UNFILTER  $\mathcal{G}$ FILTERED CORRECT DEPTH (SP) USCS CODE PROJECT: of black ranging 7 grained Jenses 4113 β. graded predominanthy mud are pyc-5,26 5, 12 333 - MWOI Ag-mg len frem BORING NO. finis MAZI Linch lenses

SOIL/SEDIMENT DESCRIPTION

LOG OF BORING

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LOC	CATIC	о ис	FBC	RIN	G:								PROJECT: CTO		BORING NO.			
													site	333	TOTAL DEPTH: 13.01			
													J08 NO.	7113	LOGGED BY: G. Helms			
													PROJ.MGR:	G. Goode	EDITED BY:			
													DRILLING CONTRACTOR;	Groundu	valer Protection			
5	00	50	sil	2	30 Y	-ir	29	احما	عامنح	n	Mo	zρ	DRILL RIG TYPE:	D130 C	(Diedrick)			
		_		-			J					Ì	DRILLER'S NAME	Charles	Bucher			
													SAMPLING METHODS:	HSA / S	plitspoon			
													STARTED TIME:	1841	DATE: 6/12/96			
													COMPLETED TIME:		DATE:			
													BORING DEPTH (fL)					
CA													CASING DEPTH (ft.)	MY VM				
Ĺ	F												WATER DEPTH (ft.)					
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NS = NO CARBON FILTERED SAMPLE READ

SOIL/SEDIMENT DESCRIPTION										
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PROJECT: NO. BORING NO.	<u>.</u>				URE	.D.	N	s		1
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LOG OF BORING

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													PROJ.MGR:	6.6000		EDITED BY:
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5	ee	-	100	1 T		* 14	ا ری	00	M	່ຼ	<b>)</b>		DRILL RIG TYPE:	Myrm)	017	20 C (Diedrick)
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													SAMPLING METHODS:	HSA /	ن مذ	it Spean
												- {	STARTED TIME:	0314		DATE: 6/13/96
												- {	COMPLETED TIME:			DATE:
	BORING DEPTH (ft.) /2.0 '															
												CASING DEPTH (fL)				
													WATER DEPTH (ft.)			
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SAMPLE DEPTH	SAMPLER TYPE	BLOWS/6-IN	NCHES DRIVEN	NCHES RECOVERED	MOISTURE	ODOR	FIL.	FILTERED OVA (PPM)	RRI	DEPTH IN FEET	1000 3031	3				
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NS = NO CARBON FILTERED SAMPLE READ

 $\vec{a}$ ū \_ DEPTH TYPE 2 BLOWS ھ DRIVEN REC'VD' 1 Z MOISTURE ODOR L. UNFILTER 8 FILTERED CORRECT DEPTH Ž Ž Ž USCS CODE PROJECT graded, Some १ व 10'-12'overall dk 1/16 11 7.0' organic organic med grains, 41.4.1 Silty SAND, fine grains Sand is a Lenses ma Peria NO harizontal are Silty and BORING NO. 333- MW03 poorly Bram ~ 6.81 color. Abundant layers

SOIL/SEDIMENT DESCRIPTION

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### APPENDIX E

## HEADSPACE METHODOLOGY FOR DETERMINING SOIL ORGANIC VAPOR CONCENTRATIONS

## HEADSPACE METHODOLOGY FOR DETERMINING SOIL ORGANIC VAPOR CONCENTRATION

Soil headspace readings where obtained utilizing the following method which conforms to the requirements of Rule 62-770.200(2), FAC.

Two 16 ounce glass soil jars were half-filled with soil sample (duplicate samples). The soil jars were then sealed utilizing "mason jar" type open top screw on caps with foil in place of the conventional solid jar tops. The soil samples were allowed to equilibrate to ambient temperature which was within the FDEP temperature range.

The samples were tested with a Foxboro Century 128, an organic vapor analyzer (OVA) equipped with a flame ionization detector (FID). Prior to each days activities, the OVA was field calibrated with 100 ppm methane in air, in accordance with the manufacturers specifications. Sample testing was performed by inserting the OVA probe through the foil sample cover and recording the highest OVA reading. Following collection of this OVA reading, the OVA was fitted with a granular activated carbon filter probe. The OVA was then used to test the headspace above the duplicate sample. Carbon absorbs petroleum hydrocarbons and thus the filtered reading is assumed to represent naturally occurring organic vapors.

Upon completion of the screening exercise, the carbon filtered result was subtracted from the unfiltered result, to obtain a net petroleum vapor value. In accordance with Rule 17(62)-770.200(2), F.A.C., and Guidelines for Assessment and Remediation of Petroleum Contaminated Soil (May 1994) corrected headspace levels in excess of 50 ppm is defined as excessively contaminated soil for diesel and used oil contaminated soil. Corrected headspace levels in excess of 10 ppm but less than 50 ppm are considered as contaminated, though not excessively contaminated.

# APPENDIX F WELL COMPLETION LOGS

Water Mgmt. Dist.:

NWFWMD

Site Information:

Permit Number:

Name:

Work Order:

6028

Address: Site G9, 323 and 333 C,S,Z: Panama City, Florida

Type of Well:

Monitor

S/T/R: Client / Consultant Information

Well Number: Method Used: Borehole Dia.

PCY333MW01 41/4" HSA 8"

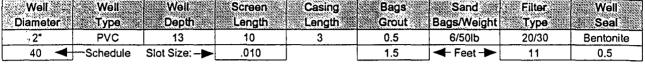
Consultant: Field Rep:

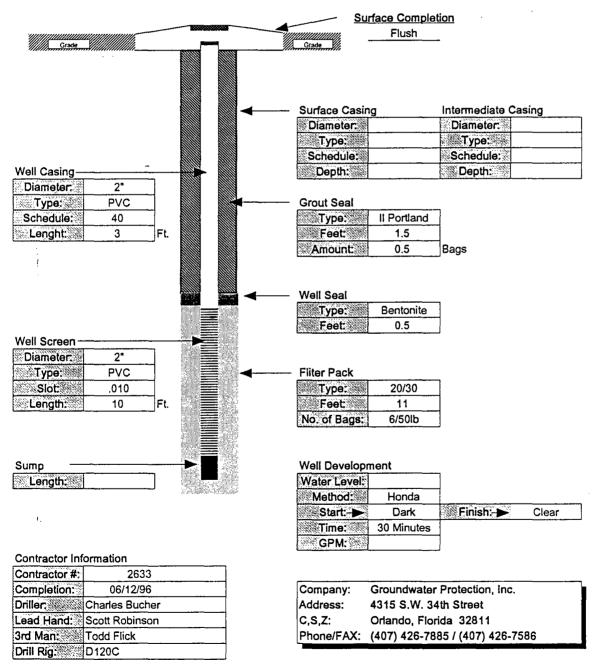
Brown & Root Gerry Goode

	Bags		114_23
] Casing	Bags	Sand F	шег

÷ , ;

CSS Panama City - Phase 1





Water Mgmt. Dist.:

**NWFWMD** 

Site Information:

Permit Number:

Name: Address C.S.Z:

: CSS Panama City - Phase 1

Address:

Site G9, 323 and 333 Panama City, Florida

Work Order. 6028

Type of Well: Monitor

S/T/R:

Type of Well: Monitor
Well Number: PCY333MW02

Client / Consultant Information

Well Number: PC:
Method Used: 4

Client / C

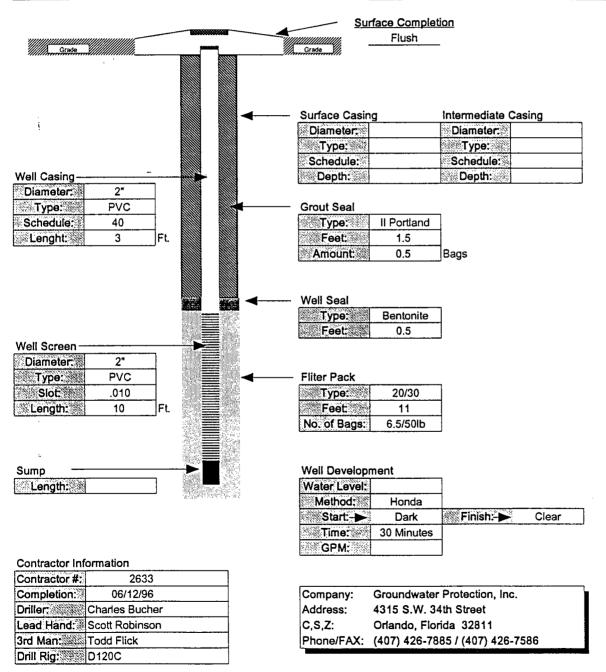
Method Used: 4½" HSA

Consultant: Brown & Root

Borehole Dia. 8"

Field Rep: Gerry Goode

Well Diameter	Well Type	Well Depth	Screen Length	Casing Length	Bags Grout	1 - 4 - 6 - 6 - 6 - 6 - 6 - 6 - 6 - 6 - 6	1100001	Well Seal
2"	PVC	13	10	3	0.5	6.5/50lb	20/30	Bentonite
40 ◀	Schedule	Slot Size:-▶	.010		1.5	<b>⋖</b> - Feet -▶	11	0.5



Water Mgmt, Dist.:

**NWFWMD** 

Site Information:

Permit Number:

Name: CSS Panama City - Phase 1

Address: Site G9, 323 and 333 Panama City, Florida

Work Order: Type of Well: 6028

C,S,Z: S/T/R:

**Monitor** 

Client / Consultant Information

Well Number: Method Used: PCY333MW03 41/4" HSA

Borehole Dia.

<u>8"</u>

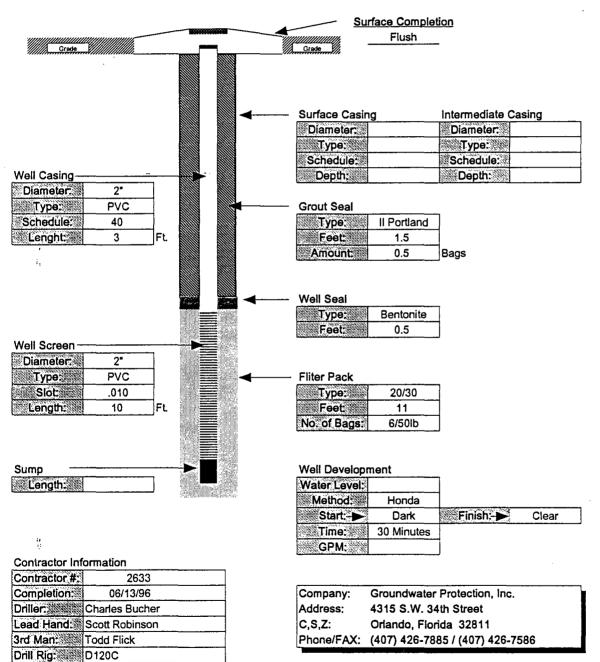
Consultant:

Brown & Root

Field Rep:

Gerry Goode

Well Diameter	Well Type	Well Depth			Bags Grout			Well Seal
2"	PVC	13	10	3	0.5	6/50lb	20/30	Bentonite
40 ◀	—Schedule	Slot Size:-	.010		1.5	Feet -►	11	0.5



Water Mgmt, Dist.:

NWFWMD

Site Information:

Permit Number:

Work Order:

Type of Well:

Method Used:

Borehole Dia.

Well Number: PCY333MW04

6028 Monitor

41/4" HSA

8"

Name: CSS Panama City - Phase 1

Address:

Site G9, 323 and 333

S . . .

C,S,Z:

Panama City, Florida

S/T/R:

Client / Consultant Information

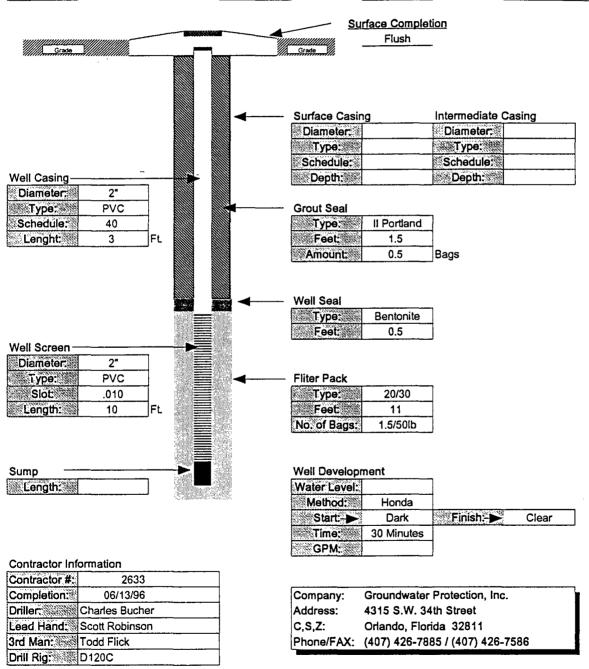
Consultant:

Brown & Root

Field Rep:

Gerry Goode

Well Diameter							Filter Type	
2*	PVC	13	10	3	0.5	1.5/50lb	20/30	Bentonite
40	Schedule	Slot Size:→	.010		1.5	<b>←</b> Feet <b>→</b>	11	0.5



### **APPENDIX G**

# GROUNDWATER GRADIENT, GROUNDWATER FLOW AND TRANSMISSIVITY CALCULATIONS

### **GROUNDWATER FLOW GRADIENT**

The groundwater flow gradient was determined using the following equation

$$i = (h_1 - h_2)/d$$

where:

i = the hydraulic gradient

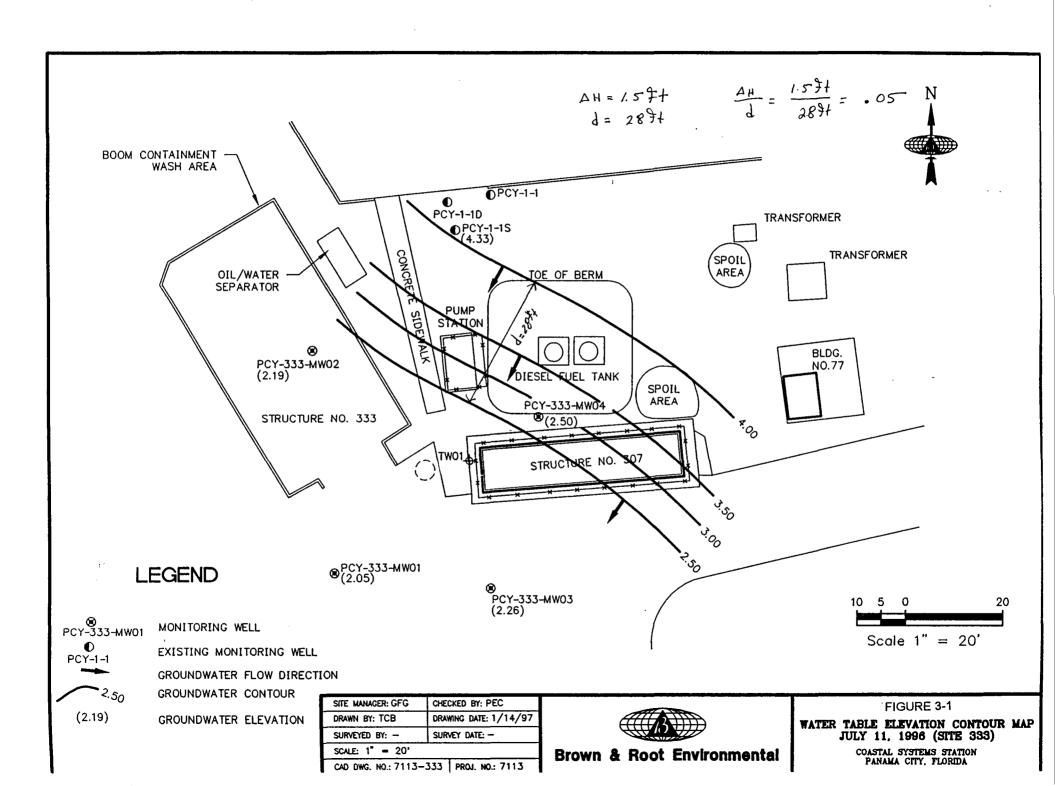
h<sub>1</sub> = the water elevation at point 1
 h<sub>2</sub> = the water elevation at point 2
 d = the distance between point 1 and point 2

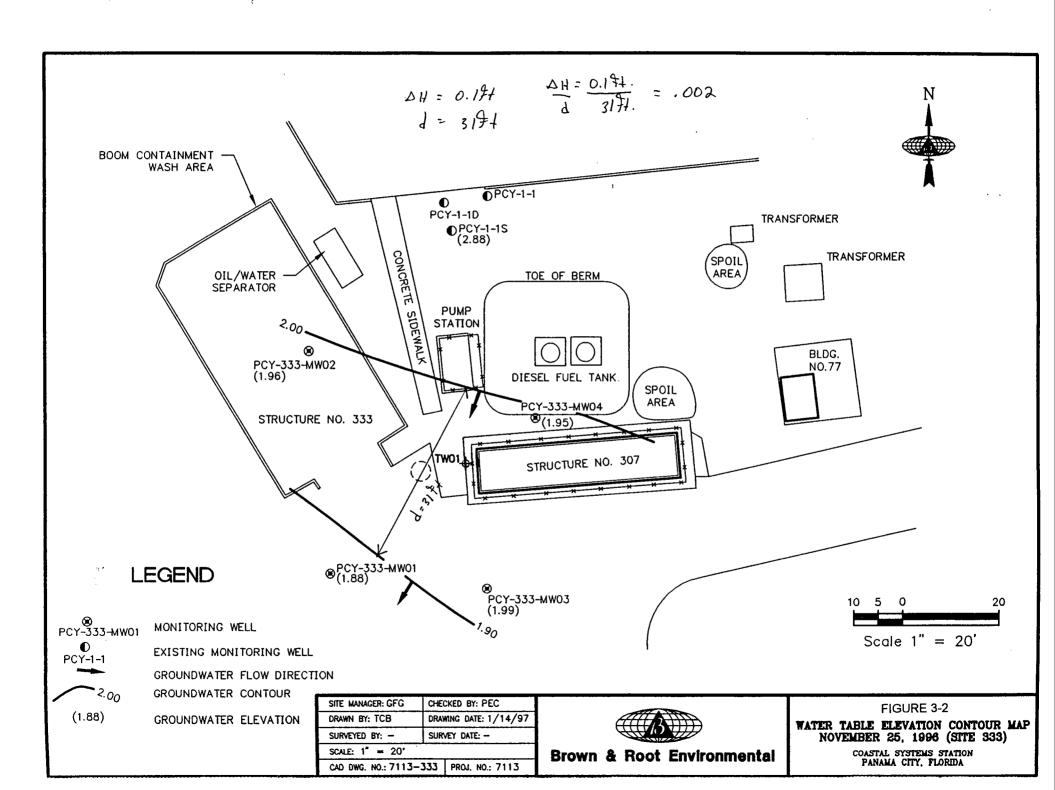
The distance and groundwater elevations were obtained from Figure 3-2.

July 11, 1996

The gradient across the site was calculated after constructing groundwater contours from the July 11, 1996, depth to water data, determining the perpendicular distance between two of these contours, and utilizing the following calculation:

$$i = 0.05 \text{ ft/ft}$$





### **GROUNDWATER FLOW VELOCITY**

Potential movement of groundwater at the site may be described in terms of transportation by natural flow system in the saturated zone while assuming groundwater flow follows Darcy's Law. Darcy's Law may be expressed as:

$$V = \frac{(K^*i)}{n_e}$$

where:

V = average velocity

K = hydraulic conductivity (0.00498 ft/min or 7.1 ft/day)

i = hydraulic gradient (0.05ft/ft)
n<sub>e</sub>= effective porosity

Therefore:

$$V = 7.1 \text{ ft/day} \times 0.05 \text{t/ft}$$

V= 1.18 ft/day

### **TRANSMISSIVITY**

Transmissivity can be determined by multiplying the hydraulic conductivity by the effective aquifer thickness ( $b_e$ ). The effective aquifer thickness is defined as depth to the top of the water table to (approximately 5 feet bls) to the top of a sandy clay unit (identified in the study area at 27 ft bls) The transmissivity was calculated as follows:

 $T = K*b_e$ 

where:

T = transmissivity

K = hydraulic conductivity (7.1 ft/day) b<sub>e</sub> = affected aquifer thickness (22 ft)

 $T = 7.1 \text{ft/day} \times 22 \text{ft}$ 

 $T = 156 \text{ ft}^2/\text{day}$ 

## APPENDIX H

## FIELD MEASUREMENTS AND SAMPLING FORMS

Subject FIELD DOCUMENTATION	Number SA-6.3	Page 17 of 32
•	Revision	Effective Date
	0	03/01/96

### ATTACHMENT C-1 EXAMPLE GROUNDWATER LEVEL MEASUREMENT SHEET

	DWATER LEV REMENT SHE	<del></del>	Page of
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PROJECT NAME: #TO-0008 PROJECT NUMBER: 713	MEASU	TION: 333	water Level Indicate
PERSONNEL: C. Buch DATE: 7-11-96 WEATHER CONDITIONS: HOTCOM	REMAR	RKS:	OR:
	WaterLevel	=Groundwater:	
Number	Reading (Feet)	Elevation: ::   Feethall	Complems
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muoz muoz		<u>4.20</u> 4.31	
mwo4		5,25	
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PC4-15		5.15	
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Signature(s)

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Cii	ent	Site	ı0:		CTO 00		Facility Address: _	'asson	1 542	=72-	76V	\		
We	ath	er:		H	OT, Char,	SLIGHTS (CENT	Sampling Method: 1	effon Bailer	ON O	her			SOP Cleaning	Y N
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## SINGLE SAMPLE LOG SHEET

Page 1 of /

Project Site Name: 10-00%	Sample I	D No.: <u>333-Gω-</u>	PC4-125-001
Project No.: 7113	Sample l	ocation: PCY-15	ss
☐ Surface Soil ☐ Subsurface Soil ☐ Sediment ☐ Other Ground which ☐ QA Sample Type:		8y: <u>C.Bursa</u>	
Sample Method:		Composite Sample E	lata
Disposable Terlan Bailer	Sample	Time	Calar/Description
Depth Sampled:			
5.15 TO 1400			
Sample Date and Time:			
7-11-96/1530			
Type of Sample			
<b>¥</b> Grab			
☐ Composite☐ Grab-Composite		Grab Sample Dat	8
☐ High Concentration	Calar	T _	Clay, Dry, Moist, Wat, etc.)
☐ Low Concentration	00.01		
Analysis Co	ntainer Requirements	Coffected (✓) Ma	n·
			ρ.
(06)/(007	40ml		
8260	40ml		,
504	125-		
presence codium chronium	540ml		
610	1172		
418.1	11.70		
8270	21.Ter		•
234.2 Observations/Notes:	500ml		
Cacle of Applicable:		Signature(s):	!
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### SINGLE SAMPLE LOG SHEET

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Project No.: 7113  Surface Soil Subsurface Soil Sediment Othe Granduster OA Sample Type:		le Location: MW03	
☐ Subsurface Soil☐ Sediment☐ Sediment☐ Street Ground uster	Samni		
☐ Sediment  ▼ Othe Ground uniter	Camp	led By: C. Bue	
14 Othe Groundwater	6.0.6	·	
	0.0.0	C. Na.:	
C CA Sample Type.	<del></del>		
Sample Method:	•	Composite Sæmp	o Data
	Samala	Time	
Disposable Texton Bailes	Sample	1 ime	Color/Desc
Depth Sampled:			
4.31 TO 12.35			
Sample Date and Time:			
7-11-96/1801			
Type of Sample			
⊠ Grab □ Composite			
☐ Grab-Composite		Grab Sample (	Data ***
☐ High Concentration	Color	Description: (San	d, Clay, Dry, Moist, We
☐ Low Concentration			
Analysis	िलाकातमः सम्वयंतम्बत	its Collected (4)	Map:
601/602	40ml		
8260	40ml		
8260	4000)		
8260 504 239,2	125-1		
8260 504	125ml 500ml		
8260 504 239.2 Accenic, cadmiun, chronin	125 ml 500 ml 500 ml		



Project No.: 7113

Sample Method:

Deoth Sampled:

☑ Grab ☐ Composite

☐ Surface Soil ☐ Subsurface Soil ☐ Sediment

1 Other Grandwiles ☐ QA Sample Type:

Disposable Tation Bailer

4.20 70 13.20 Sample Date and Time:

7-11-96 / 160Z

☐ Grab-Composite ☐ High Concentration

☐ Low Concentration

Analysis ...

601/602

42G0 504

41811

8270 239.2

Type of Sample

SINGLE SAMPLE LOG SHEET Page Sample ID No.: 333-GW-mwoz-001 Project Site Name: 270-000x 333 Sample Location: \_\_\_\_\_oz Sampled By: C. Burg. C.O.C. No.: \_\_\_\_\_ Composite Sample Data Sample Time Calar/Description Grab Sample Data Calar Description: (Sand, Clay, Dry, Moist, Wet, etc.) Container Requirements Collected (V) Map: Yoml ·40~1 125ml Liec 1 Liter 50ml

	adin chemital	50ml		<del></del>	
Observation	ins/Notes:				
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MS/MSD | Duplicate ID No:

### SINGLE SAMPLE LOG SHEET

Sample ID No .: 333-6W-mwo 1-001 Project Site Name: CTS-0008 333 Sample Location: mwol Project No.: 7113 Sampled By: C. Burgan ☐ Surface Soil ☐ Subsurface Soil ☐ Sediment C.O.C. No.: \_\_\_\_\_ 1 Other Grandwiter ☐ QA Sample Type: Sample Method: Composite Sample Data Disposable to la Ruiler Depth Sampled: Sample Time Calar/Description 4.60 To 13.20 Sample Date and Time: 7-11-96/1714 Type of Sample **⊠** Grab ☐ Composite Grab Sample Data ☐ Grab-Composite ☐ High Concentration Calar Description: (Sand, Clay, Dry, Moist, Wat, etc.) ☐ Low Concentration Collected (4/) Analysis... Container Requirements 40 -1 601/602 40ml 8260 125-1 504 418.1 11,700 610 11ites 2312 500 ml 8270 2 Liter Assonic Cadmium Chromium 5000 Observations/Notes: Carcle if Applicable:

Brown & Root Environmental							SAMPLING LOG Page / of /								
888	2.0	ect	No:	5	7113		Date: 7-11-96 Sampler(s): C, Bus un								
Ciie	nt S	ite !	o: <u>(</u>	<u> </u>	00088		Facility Address:	<u> محتر</u>	1395	EM STR	Tre	m Re	P		
We	the	r: H	61	cl	20,51.36	TBreze	Sampling Method: 1	etlon Bailer	<b>⊘</b> N Oth	ner			SOP Cleaning	YN	
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GRORDOWATER SYSTEM PERFORMANCE AND QUALITY CONTROL SAMPLES TEST PARAMETERS															
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Subject FIELD DOCUMENTATION	Number SA-6.3	Page 17 of 32
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	0	03/01/96
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### ATTACHMENT C-1 EXAMPLE GROUNDWATER LEVEL MEASUREMENT SHEET

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	3			WATER LE				
			MEASUR	KEMEN 3H	<u> </u>	Page I of		
	PROJECT NAI PROJECT NU PERSONNEL: DATE: 11/0 WEATHER CO	MBER:	7113	MEASI ADJU	TION: <u>5/7E</u> JRING DEVICE STMENT FACT RKS: <u>M/1144</u>	E: <u>Johnso Probe</u> TOR:		
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			Reference Point	indicator Reading: IFeeti C	Elevation: (Feet)	Companis		
	PCY-15 -	10:15		6.60		Total Raph 15. 20'		
	MW03	10:20		4.58		Jaho / Repll 9.45		
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	MWOZ	10:40		4.40		Total Repth 13.10		
	TWOI	10:50	·	6.76		THE Rept 10.23		
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Brown & Root Environmental BRE Project No: 7/13							SAMPLING EOG   Page / of /     Date: 11/25/96   Sampler(s): 6(-a/d 60000								
Client Site ID: 5:14 333							Facility Address: (001h) System Shaffor								
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					113		Date: 11/25/96   Sampler(s): 6-3/ 6-6								
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Project Site Name: Could	Systems .	Show Sampi	e ID No.: <u>333</u> - 6	6W-72	01-001B
Project No.: 71/3			e Location:	ank	
☐ Surface Soil ☐ Subsurface Soil ☐ Sediment ☐ Other <u>Black</u> (9100) ☐ QA Sample Type:	Sue Les)		ed By: <u>Garald</u> . No.: <u>009</u> -		,
Samole Method: Bailor/ Prishellic Prof			Composite San		Calado
		Sample	Time	<del></del>	Color/Description
Depth Sampled:	<u> </u>				Clear
- N/A		<u> </u>	<u> </u>		
Sample Date and Time:					
11/25/96 / 12:15					
Type of Sample					
☐ Grab ☐ Composite					
☐ Grab-Composite			Grab Sample	a Data	
☐ High Concentration	C	olor	Description: (s	and, Clay, O	ry, Maiet, Wat, atc.)
☐ Low Concentration					
Апаузіз	Contair	ве Каригалал	ts Collected (4)	Мар:	
601	40 ml	usal, gloss			
602	i	vial, 1/11			
8260	1	vial, g/iss			
610	<b>.</b>	en hun, 9/011		1	/A
8270	1	b , 1/01	, /	7	-
504	ì	. 1''		1	
Metal, Accenic Codnim Ch.	1	,		1	
418,1 (TRPH)	1603	· 1	,	1	
Observations/Notes:	1		· · · · · · · · · · · · · · · · · · ·		
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Project Site Name: Conhol	Zileni Strian	Sample IC	No.: 333-6	SW - MWO4 - 0	00/
Project No.: 7113		Sample L	ocation:	or well in	104
☐ Surface Soil☐ Subsurface Soil☐		Sampled	By: <u>Genl</u>	1 600de	
☐ Sediment		C.O.C. N	o.: <u>00 9</u> 2	7	
☑ Other <u>Grown Auctor</u> ☐ QA Sample Type:	· · · · · · · · · · · · · · · · · · ·				
Sample Method:			Composite Sam	pia Data	
Bailer/Peristaldic P.	<u></u>	nple	Time	Calar/0	Description
Depth Sampled: with Tall	<u></u>			Chor/o	ily
monitor well					<i>,</i>
Sample Date and Time:					
11/25/96 / 14:45					
Tvoe of Samole  Grab					
☐ Composite					
☐ Grab-Composite			Grab Sample	· · · · · · · · · · · · · · · · · · ·	
☐ High Concentration☐ Low Concentration	Color		Description: (s.	and, Clay, Ory, Mais	t, Wat, atc.)
		~		1	<del></del> -
Апаузіз	Container Requ	arements	Collected (A)	Map:	
601	Youl vist	5/111		. 90	4-15
602	40 2/ 41)	9/111		-	
8260	40 ml vial	gless		ح د د د د	* MWOY
610	1 / honder,	Glist			Twal
8270	26-6-1	191011			
504	125 ml and	~ · / · · · · · · · · · · · · · · · · ·		1	
metals, Lead, aigmic, Cd, Chi.	1 ' 1	1		Ano!	3 com
Observations/Notes: Page	Free produce		15' thic	knoss From w	ele
prior to Sompling		/		·	
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Project Site Name: Coashal	Sychm Shelsa	· Sample I	D No.: 333-61	w-Twol	- 001
Project No.: 7113		Sample t	ocation:	b, wel	Two I sil 333
☐ Surface Soil☐ Subsurface Soil☐		Sampled	By: Gorald (	Gooda	
☐ Sediment ☐ Other <u>Groudus</u> — ☐ QA Sample Type:		C.O.C. N	la.: <u>0097</u>		
Sample Method: Bile(/			Composite Sam	pie Data	
Peristollic purp	5	Sample	Time		Calar/Description
Depth Sampled: water Toble  — monitor well	, -	<del>-</del>			Cleal
Sample Date and Time:					
Type of Sample  ☐ Grab ☐ Composite ☐ Grab-Composite			Grab Sample		
☐ High Concentration ☐ Low Concentration	Color		Description: (Sa	and, Clay, D	ry, Maist, Wat, etc.)
Analysis 66 /	Container R		Collected (4)	Мар:	· pcy -15
60Z 8260	40 ml vial 40 ml vial 40 ml vial	13, 9/201		• Awoz	Young
610 8270	1 hombos	1, 9/015			- 7wol
Molals Associa, Lood, Cd	Ch. / Lp1.			• 4400	l . 
Observations/Notes:	1- 110512	1607 mln;	(1/6/11)	<u></u>	·
MS/MSD Ouplicate ID No:	Collected of		Signature(s):	. Tood	7



Sample ID No.: 333 - 6w - Mwo2 - 002 Project Site Name: Cooked Systems Station Sample Location: Monito, well There? Project No.: 7113 Sampled By: Gud Horde ☐ Surface Soil ☐ Subsurface Soil C.O.C. No.: 0097 ☐ Sediment E Other Granduste ☐ QA Sample Type: \_ Composite Sampla Data Sample Method: Penistaltic Porp Calar/Description Sample Time Depth Sampled: waln 7.36 Clan minitor well Sample Date and Time: 11/25/96 / /2:10 Type of Sample ☑ Grab ☐ Composite Grab Sample Data ☐ Grab-Composite ☐ High Concentration Description: (Sand. Clay, Ory, Maist, Wat, atc.) Calar ☐ Low Concentration Container Requirements Collected (4) - zizyłanA Pcy-15 1 h plante Observations/Notes: Signature(s): Carcle of Applicable: New Yords MS/MSD Ouplicate ID No:

# APPENDIX I

# **GROUNDWATER FIELD SCREENING RESULTS**

Job: "711	.3BTEX" (Standard A, )	06/12/96	09:54:11		
Petector "A" ( gment Width - 10ise - 0 Detector Temp - 110 C	Attenuation - Range - Baseline - Oven Temp -	1 1 70 70 C	Min Area Autorange Units	-	0 1 PPM
Peak Component Name	Concentration	Height	Area	Time	Alarm
Detector "B" ( PID Segment Width - 10 Noise - 10 Detector Temp - 110 C	Attenuation - Range - Baseline - Oven Temp -	10 1 435 70 C	Min Area Autorange Units	<u>-</u>	0 1 PPB
Peak Component Name	Concentration	Height	Area	Time	Alarm
1 *** unknown ** 2 BENZENE 3 TOLUENE 4 ETHYLBENZENE 5 M, P-XYLENES 6 O-XYLENES	* 65.278 100.000 100.000 100.000 100.000	995 18847 23272 17408 47235 17802	230570 2776070 4290270 2849640 11732860 4807930	1:17 1:40 2:26 3:54 4:05 4:39	
-	End of Result	Report			

Job "7113BTEX"	' (333-GW-TW03-003	L) 06/12/9	6 11:53:18		
Detector "A" ( ) Sigment Width - 10 ise - 0 Letector Temp - 110 C	Attenuation - Range - Baseline - Oven Temp -	1 1 0 70 C	Min Area Autorange Units	- - -	0 1 PPM
Peak Component Name	Concentration	Height	Area	Time	Alarm
Detector "B" ( PID) Segment Width - 10 Noise - 5 Detector Temp - 110 C	Attenuation - Range - Baseline - Oven Temp -	10 1 472 70 C	Min Area Autorange Units	- -	0 1 PPB
Peak Component Name	Concentration	Height	Area	Time	Alarm
1 *** unknown ***		1617	359050	1:10	
	End of Result	Poport			

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Job "7113BTEX"	(333-GW-TW04-001	.) 06/12/9	6 12:12:32		
Detector "A" ( ) Sagment Width - 10 ise - 0 Detector Temp - 110 C	Attenuation - Range - Baseline - Oven Temp -	1 1 0 70 C	Min Area Autorange Units	-	0 1 PPM
Peak Component Name	Concentration	Height	Area	Time	Alarm
Detector "B" ( PID) Segment Width - 10 Noise - 5 Detector Temp - 110 C Peak Component Name	Attenuation - Range - Baseline - Oven Temp - Concentration	10 1 450 70 C Height	Min Area Autorange Units Area	Time	0 PPB Alarm
1 *** unknown *** 2 *** unknown *** 3 *** unknown *** 4 *** unknown ***		2404 51 42 636	389360 3760 4370 296710	1:11 2:19 7:48 9:34	
	End of Result	Report			

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Job "7113BTEX"	(333-GW-TW05-001	L) 06/12/9	6 12:30:02		
Detector "A" ( ) Segment Width - 10 ise - 0 Letector Temp - 110 C	Attenuation - Range - Baseline - Oven Temp -	1 1 0 70 C	Min Area Autorange Units	-	0 1 PPM
Peak Component Name	Concentration	Height	Area	Time	Alarm
Detector "B" ( PID) Segment Width - 10 Noise - 6 Detector Temp - 110 C	Attenuation - Range - Baseline - Oven Temp -	10 11 448 70 C	Min Area Autorange Units	- - Time	0 1 PPB
Peak Component Name	Concentration	Height ———	Area	1 I I III e	Alarm
1 *** unknown ***		493	75460	1:02	
	End of Result	Report			

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	Job ["7113BTEX"	(333-GW-TW06-001	.) 06/12/9	6 12:45:53		
ıse	tor "A" ( ) nt Width - 10	Attenuation - Range - Baseline - Oven Temp -	1 1 0 70 C	Min Area Autorange Units	-	0 1 PPM
Peak	Component Name			Area	Time	Alarm
Detec Segme: Noise Detec	tor "B" ( PID) nt Width - 10 tor Temp - 110 C	Attenuation - Range - Baseline - Oven Temp -	10 1 459 70 C	Min Area Autorange Units	- - -	0 1 PPB
Peak	Component Name			Area	Time	Alarm
123456789012345678901234567	*** unknown ***  *** unknown ***  *** unknown ***  BENZENE  *** unknown ***  TOLUENE  TOLUENE  *** unknown ***  *** unknown ***  ETHYLBENZENE  M,P-XYLENES  *** unknown ***  O-XYLENES  *** unknown ***  *** unknown ***  *** unknown ***  *** unknown ***  *** unknown ***  *** unknown ***  *** unknown ***  *** unknown ***  *** unknown ***  *** unknown ***  *** unknown ***  *** unknown ***  unknown ***  unknown ***  *** unknown ***  *** unknown ***  *** unknown ***  *** unknown ***  *** unknown ***  *** unknown ***  *** unknown ***	14.383 19.473 23.685 136.917 43.166 169.125	002023564550890518911111142 362859512225201497942228266 230406720229784692493067183 111244689678806323544644532	244440718600000000000000000000000000000000000	71706213161723471384435469202244023525012303500240241555151111112222233444455556667778889	

End of Result Report

Job "7113BTEX"	(Standard B, )	06/12/96	10:08:51		
Detector "A" ( )  "gment Width - 10  ise - 0  Detector Temp - 110 C	Attenuation - Range - Baseline - Oven Temp -	1 1 0 70 C	Min Area Autorange Units	- - -	0 1 PPM
Peak Component Name	Concentration	Height	Area	Time	Alarm
Detector "B" ( PID) Segment Width - 10 Noise - 7 Detector Temp - 110 C	Attenuation - Range - Baseline - Oven Temp -	10 1 444 70 C	Min Area Autorange Units	<u>-</u>	0 1 PPB
Peak Component Name	Concentration	Height	Area	Time	Alarm
1 *** unknown *** 2 *** unknown *** 3 BENZENE 4 TOLUENE 5 ETHYLBENZENE 6 M,P-XYLENES 7 O-XYLENES	17.168 23.591 17.729 18.681 17.749 End of Result	19842 984261 2733901 262621 2827	8960 223150 403500 1070260 411150 2074860 686380	1::427 1::427 4::509	

7.1: U.7.1.2.D.T.T.V.	/G 1 1 G 1	06/10/06	10 22 10	<del></del>	
Job "7113BTEX"	(Standard C, )	06/12/96	10:22:40		
Detector "A" ( ) Sigment Width - 10 Lise - 0 Detector Temp - 110 C	Attenuation - Range - Baseline - Oven Temp -	1 0 70 C	Min Area Autorange Units	- - -	0 1 PPM
Peak Component Name	Concentration	Height	Area	Time	Alarm
Detector "B" ( PID) Segment Width - 10 Noise - 4 Detector Temp - 110 C	Attenuation - Range - Baseline - Oven Temp -	10 432 70 C	Min Area Autorange Units	- - -	0 1 PPB
Peak Component Name	Concentration	Height	Area	Time	Alarm
1 *** unknown *** 2 BENZENE 3 TOLUENE 4 ETHYLBENZENE 5 M,P-XYLENES 6 O-XYLENES	6.182 12.116 3.9961 5.046	734 1066 4203 766 2617 788	192970 147720 660110 109720 606770 178930	1:12 1:38 2:23 3:48 3:59 4:31	
0 O-VIDENES	5.010				

\$ ,1

### **APPENDIX J**

# **GROUNDWATER LABORATORY DATA SHEETS**

(Groundwater Samples Collected July 11 and November 25, 1996)

# **GROUNDWATER SAMPLES COLLECTED JULY 11, 1996**

# CASE NARRATIVE GC/MS VOLATILE ORGANICS

QAL Lab	Reference	No./SDG.	MB:	370	
Project	Brow	n & Root	Coastal	Systems	Station

#### I. RECEIPT

No exceptions were encountered unless a Sample Receipt Exception Report is attached to the Chain-of-Custody included with this data package.

#### II. HOLDING TIMES

- A. Sample Preparation: N/A
- B. Sample Analysis: All holding times were met.

#### III. METHOD

Preparation: N/A Cleanup: N/A

Analysis: SW-846 8260

#### IV. PREPARATION

Not applicable.

#### V. ANALYSIS

- A. Calibration: All acceptance criteria were met.
- B. Blanks: All acceptance criteria were met.
- C. Surrogates: All acceptance criteria were met.
- D. Spikes: As requested, the matrix spikes were performed using a sample from sample delivery group MB370 (MB370003MS and MB370003MSD). Please note that the relative percent recovery for 2-Chloroethylvinyl ether was above the advisory QC criteria. A copy of the results is provided for your review.
- E. Samples: Sample analysis proceeded normally.
- F. Other: Please note that the Form 1's reflect the specified target list.

Currently, there are not enough data points collected to produce control charts for the water surrogate recoveries for 1,2-Dichloroethane-d4. These charts are in the process of being developed.

Lab Reference No./SDG: MB370
Page 2

A summary of the most current applicable method detection limits (MDLs) immediately follows the case narrative.

I certify that this data package is in compliance with the terms and conditions agreed to by the client and QAL, Inc., both technically and for completeness except for the conditions noted above. Release of the data contained in this hardcopy data package has been authorized by the Laboratory Manager or designated person, as verified by the following signature.

SIGNED:

Debra A. Vergin

Chemisť

Lab Reference No./SDG: MB370

# CASE NARRATIVE Addendum

Sample Information

LAB SAMPLE ID	CLIENT SAMPLE ID	SAMPLE MATRIX	DATE SAMPLED	DATE EXTRACTED	DATE <u>ANALYZED</u>	SAMPLE pH <sup>1</sup>
MB370001	333-MW01-1	WATER	07/11/96	N/A	07/18/96	<2
MB370001	333-MW02-1	WATER	07/11/96	N/A N/A	07/18/96	<2
MB370003	333-MW03-1	WATER	07/11/96	N/A	07/19/96	<2
MB370003MS	333-MW03-1MS	WATER	07/11/96	N/A	07/19/96	<2
MB370003MSD	333-MW03-1MD	WATER	07/11/96	N/A	07/19/96	<2
MB370004	333-PC4-1	WATER	07/11/96	N/A	07/19/96	<2
MB370005	333-MW01-1B	WATER	07/11/96	N/A	07/19/96	<2
MB370006	333-MW03-1D	WATER	07/11/96	N/A	07/19/96	<2
MB3700.07	TRIP_BLANK	WATER	07/11/96	N/A	07/18/96	<2
X07186B1	VBLKOS	WATER	N/A	N/A	07/18/96	N/A
X07196B1	VBLK0V	WATER	N/A	N/A	07/19/96	N/A

<sup>1</sup> Applies to samples designated for purgeable VOA analysis only.

#### ORGANICS ANALYSIS METHOD DETECTION LIMITS

Laboratory	Name:	CH2M HILL	Sample Matrix:	WATER
Analytical	Method:	SW8260	Extraction Method:	
			10	7.T
			-	OL.
	3 t		<u>ug</u>	
	Acet	one conitrile	1.7	
		olein	7 5.:	
		rlonitrile	4.9	
	Benz		0.4	•
		obenzene	0.:	
		ochloromethane	0.:	
•		odichloromethane	0.:	
		ofluorobenzene	1.4	
•		oform	0.4	
		nomethane	0.9	
		itanone	0.	
		on Disulfide	0.4	
		on Tetrachloride	0.:	
•		robenzene	0.9	
		roethane	0.:	
		loroethylvinyl eth		
		roform	0.:	
		romethane	0.:	
		proprene	0.1	
		loropropene	0.:	
		lorotoluene	1.:	
		lorotoluene	0.1	
	· ·	1,2-Dichloroethene		
		1,3-Dichloropropen		
		ohexanone	33.4	
	-	Dibromo-3-chloropr		
		comochloromethane	0.:	
	1,2-	Dibromoethane	0.:	
	Dibr	comofluoromethane	1.9	97
	Dibr	comomethane	0	42
· ·	1,2-	Dichlorobenzene	0.	
	1,3-	Dichlorobenzene	0.	
		Dichlorobenzene	0.	
	Dich	lorodifluoromethan	ue 0.	62
	1,1-	Dichloroethane	0	42
	1,2-	Dichloroethane	0.1	
	1,1-	Dichloroethene	0.	37
	1,2-	Dichloroethene (to		
	Dich	lorofluoromethane	0.:	
	1,2-	Dichloropropane	0	
	1,3-	Dichloropropane	0	
	2,2-	Dichloropropane	0	23

#### ORGANICS ANALYSIS METHOD DETECTION LIMITS

Laboratory Name: CH2M HILL	Sample Matrix: WATER
Analytical Method: SW8260	Extraction Method:
	MDL
	ug/L
1,1-Dichloropropene	0.45
1,4-Dioxane	12.96
Ethyl ether	0.36 0.51
Ethyl methacrylate	
Ethylbenzene	0.62 1.17
Hexachlorobutadiene	0.80
2-Hexanone Iodomethane	1.47
	11.96
Isobutyl alcohol Isopropylbenzene	0.45
m-,p-Xylene	1.07
Methacrylonitrile	0.37
Methyl methacrylate	0.40
- Methyl tert-butyl eth	
4-Methyl-2-pentanone	0.39
Methylene Chloride	0.38
n-Butylbenzene	0.64
n-Propylbenzene	0.53
Naphthalene	2.30
o-Xylene	0.59
p-Isopropyltoluene	0.68
Pentachloroethane	0.31
Propionitrile	4.87
sec-Butylbenzene	0.51
Styrene	0.55
tert-Butylbenzene	0.46
1,1,1,2-Tetrachloroet	nane 0.33
1,1,2,2-Tetrachloroet	nane 0.78 ,
Tetrachloroethene	0.46
Tetrahydrofuran	5.47
Toluene	0.45
Toluene-d8	1.26
trans-1,2-Dichloroeth	ene 0.26 .
trans-1,3-Dichloropro	pene 0.43
trans-1,4-Dichloro-2-	
1,1,2-Trichloro-1,2,2	-trifluoroet 1.87
1,2,3-Trichlorobenzen	
1,2,4-Trichlorobenzen	
1,1,1-Trichloroethane	
1,1,2-Trichloroethane	0.44
Trichloroethene	0.44
Trichlorofluoromethan	
1,2,3-Trichloropropan	e 0.87

#### ORGANICS ANALYSIS METHOD DETECTION LIMITS

Laboratory	Name:	CH2M HILL	Sample	Matrix:		WATER
Analytical	Method:	SW8260	Extract	cion Met	:hod:	
					MD	L
-					ug/	L
	1,2,	4-Trimethylbenzene			0.4	4
	1,3,	5-Trimethylbenzene			0.6	5
•	Viny:	l acetate			1.1	6
	Viny	l Chloride			0.3	7
	Xyle	ne (total)			1.5	9

333-MW01-1

Lab Name: CH2M HILL

Contract: MB370

SAS No.: SDG No.: MB370

Matrix: (soil/water) WATER

Lab Sample ID: MB370001

Sample wt/vol:

Lab Code: MGM

5.0 (g/mL) ML

Case No.: MB370

Lab File ID: 18JUL0901009.D

Level: (low/med)

ed) LOW

Date Received: 07/13/96

% Moisture: not dec.

Date Analyzed: 07/18/96

Column: (pack/cap) CAP

Dilution Factor: 1.0

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Dilucion Factor: 1.

CONCENTRATION UNITS:

CAS NO.

COMPOUND

(ug/L or ug/Kg) UG/L

Q

74-87-3Chloromethane	10	ប
75-01-4Vinyl chloride	10	บ
74-83-9Bromomethane	10	Ü
75-00-3Chloroethane	10	ט
75-69-4Trichlorofluoromethane		ט
75-35-41,1-Dichloroethene		_
75-09-2Methylene chloride	10	บ บ
156-60-5trans-1,2-Dichloroethe		Ü
75-34-31,1-Dichloroethane		
75-34-31, 1-DICHIOFOECHARE	10	Ŭ
67-66-3Chloroform		Ū
71-55-61,1,1-Trichloroethane		Ū
56-23-5Carbon tetrachloride		Ū
71-43-2Benzene	10	ָט
107-06-21,2-Dichloroethane	10	Ū
79-01-6Trichloroethene		Ŭ
78-87-51,2-Dichloropropane		Ŭ
75-27-4Bromodichloromethane		Ū
110-75-82-Chloroethylvinyl eth	ner  10	Ū
10061-01-5cis-1,3-Dichloroproper	ne	Ū
108-88-3Toluene		U
10061-02-6trans-1,3-Dichloroprop	pene10	U
79-00-51,1,2-Trichloroethane		ַ ט
127-18-4Tetrachloroethene	10	ן ט
124-48-1Dibromochloromethane	10	ע
108-90-7Chlorobenzene	10	ប
100-41-4Ethylbenzene	4	J
75-25-2Bromoform	10	U
79-34-51,1,2,2-Tetrachloroeth	nane 10	ָט
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#### 1E VOLATILE ORGANICS ANALYSIS DATA SHEET TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

333-MW01-1

Lab Name: CH2M HILL

Contract: MB370

Lab Code: MGM Case No.: MB370 SAS No.:

SDG No.: MB370

Matrix: (soil/water) WATER

Lab Sample ID: MB370001

Sample wt/vol: 5.0

(q/mL) ML

Lab File ID: 18JUL0901009.D

Level: (low/med) LOW

Date Received: 07/13/96

% Moisture: not dec.

Data Analyzed: 07/18/96

Column: (pack/cap) CAP

Dilution Factor: 1.0

Number TICs found: 4

CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
CAS NUMBER  ===================================	COMPOUND NAME  Silane, methoxytrimethyl- Benzene, trimethyl- isomer Benzene, propenyl- isomer Naphthalene	RT  9.004 26.375 27.969 32.110	EST. CONC.  10 7 7 30	Q NJ NJ NJ NJ
28. 29. 30.				

SW846-8260

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Data File: /chem/ms5.i/a071896a.b/18Jul0901009.d

Report Date: 29-Jul-1996 15:29

# OAL, Inc.

VOLATILE REPORT METHOD 8260/5-ML PURGE Data file : /chem/ms5.i/a071896a.b/18Jul0901009.d/

Lab Smp Id: MB370001 Client Smp ID: 333-MW01-1

Autotune Date: 18-Jul-96 11:10

Inj Date : 18-JUL-1996 19:32 Operator : WLH/RLW Smp Info : 333-MW01-1 MB370001 Inst ID: ms5.i

Misc Info : 5MLS

Comment

: /chem/ms5.i/a071896a.b/8260w5.m Method

Quant Type: ISTD Meth Date : 29-Jul-1996 15:28 ms5

Cal File: 10Jul2001020.d Cal Date : 10-JUL-1996 22:29

Als bottle: 9

Dil Factor: 1.000 Integrator: HP RTE

Compound Sublist: BROWN&ROOT.sub

Target Version: 3.10

	-	QUANT SIG				ON-COLUMN	FINAL
Compou	inds .	MASS	RT	EXP RT REL RT	RESPONSE	( ug/L)	( ug/L)
	· · ·	***	==	*****	******	3228384	======
• 1	Pentafluorobenzene	168.00	13.755	13.756 (1.000)	1709925	50	
3	Chloromethane	50.00	Con	npound Not Detect	ed.		
4	Vinyl chloride	62.00	Соп	mpound Not Detect	ed.		
5	Bromomethane	94.00	Con	mpound Not Detect	ed.		
. 6	Chloroethane	64.00	Соп	mpound Not Detect	ed.		
7	Trichlorofluoromethane	101.00	Con	mpound Not Detect	ed.		
11	1,1-Dichloroethene	96.00	Con	mpound Not Detect	ed.		
16	Methylene chloride	84.00	Con	mpound Not Detect	ed.		
19	trang-1,2-Dichloroethene	96.00	Con	mpound Not Detect	ed.		
21	1,1-Dichloroethane	63.00	Соп	mpound Not Detect	ed.		
31	Chloroform	83.00	Соп	mpound Not Detect	ed.		
33	1,1,1-Trichloroethane	97.00	Con	mpound Not Detect	ed.		
35	Carbon tetrachloride	117.00	Con	mpound Not Detect	ed.		
* 36	1,4-Difluorobenzene	114.00	16.047	16.038 (1.000)	2186995	50	
37	1,2-Dichloroethane	62.00	Con	mpound Not Detect	ed.		
38	Benzene	78.00	Con	mpound Not Detect	ed.		
39	Trichloroethene	95.00	Con	mpound Not Detect	ed.		
41	1,2-Dichloropropane	63.00	Con	mpound Not Detect	ed.		
43	Bromodichloromethane	83.00	Con	mpound Not Detect	ed.		
45	2-Chloroethylvinyl ether	63.00	Con	mpound Not Detect	ed.		
46	cis-1,3-Dichloropropene	75.00	Con	mpound Not Detect	ed.		-
* 48	Chlorobenzene-d5	117.00	22.382	22.372 (1.000)	1678156	50	
49	Toluene	91.00	Cor	mpound Not Detect	ed.		
51	trans-1,3-Dichloropropene	75.00	Сот	mpound Not Detect	ed.		
52	1,1,2-Trichloroethane	83.00	Cor	mpound Not Detect	ed.		
55	Tetrachloroethene	166.00	Cor	mpound Not Detect	ed.		
56	Dibromochloromethane	129.00	Cor	mpound Not Detect	ed.		
58	Chlorobenzene	112.00	Cor	mpound Not Detect	ed.		
60	Ethylbenzene	91.00	22.618	22.608 (1.011)	203237	4	4 (a)
66	Bromoform	173.00	Cor	mpound Not Detect	ed.		

Page 1

A. 3

CONCENTRATIONS

								CONCENTRA	ATIONS
			QUANT SIG					ON-COLUMN	FINAL
Co	ompou	ınds	MASS	RT	EXP RT	REL RT	RESPONSE	( ug/L)	( ug/L)
	51	trans-1,3-Dichloropropene	75.00		19.874	Com	pound Not De	etected.	
	52	1,1,2-Trichloroethane	83.00		20.169	Соп	pound Not De	tected.	
	55	Tetrachloroethene	166.00		20.779	Com	pound Not De	etected.	
	56	Dibromochloromethane	129.00		21.162	Com	pound Not De	etected.	
	58	Chlorobenzene	112.00		22.461	Com	pound Not De	etected.	
	60	Ethylbenzene	91.00	22.618	22.608	(1.011)	203237	4	4 (a)
	66	Bromoform ·	173.00		24.330	Com	pound Not De	etected.	
	68	1,1,2,2-Tetrachloroethane	83.00		24.753	Com	pound Not De	tected.	
*	69	1,4-Dichlorobenzene-d4	152.00	27.329	27.320	(1.000)	856526	50	
\$	91	Dibromofluoromethane	113.00	13.775	13.766	(1.001)	924140	52	52
\$	92	1,2-Dichloroethane-d4	65.00	15.024	15.005	(0.936)	374166	44	44
\$	93	Toluene-d8	98.00	19.283	19.274	(1.202)	1937159	48	48
\$	94	Bromofluorobenzene	174.00	24.909	24.890	(1.113)	785452	47	47

# QC Flag Legend

Report Date: 29-Jul-1996 16:10

#### Quality Analytical Laboratories

Unknown Compounds Quantitation Report

Data file : /chem/ms5.i/a071896a.b/18Jul0901009.d

Lab Smp Id: MB370001 Client Smp ID: 333-MW01-1

Inj Date : 18-JUL-1996 19:32 Autotune Date: 18-Jul-96 11:10:3

Operator : WLH/RLW Inst ID: ms5.i

Smp Info : 333-MW01-1 MB370001 Misc Info : 5MLS

Comment

Method : /chem/ms5.i/a071896a.b/8260w5.m

Meth Date : 29-Jul-1996 16:09 whall

Cal Date : 10-JUL-1996 22:29 Cal File: 10Jul2001020.d

Als bottle: 9

Dil Factor: 1.000 Target Version: 3.12

Integrator: HP RTE Compound Sublist: BROWN&ROOT.sub

Sample Matrix: WATER

Quantitative Mode : Use RF of Nearest Std

Concentration Formula: Uf \* 5/Vo

Name	Value	Description
Uf Vo		ng unit correction factor Sample Volume purged (mL)

IS'	TD		RT	AREA	AMOUNT
==	===:	===	====	=====	=====
1	1	Pentafluorobenzene	13.755	7126628	50.000
*	69	1,4-Dichlorobenzene-d4	27.329	6312538	50.000

		CONCENT	RATIONS		QU	IANT	
RT	AREA	ON-COL( ug/L)	FINAL( ug/L)	QUAL	LIBRARY	LIB ENTRY	CPND #
****	***	*****	***********	****	******	********	
Silane, n	methoxytri	methyl-		CAS	3 #: 1825-61-2		
9.004	1510295	10	10	91	NBS75K.l	63613	1
Benzene,	trimethyl	isomer		CAS	5 #: <del>-526-73-8</del>	- prop 8/	5-/95-
26.375	931914	7	7	97	NBS75K.1	64573	69
Benzene,	propenyl	isomer		CAS	\$ #: <del>-873-66-5</del>	- mm 8/	5/95
27.969	851357	7 .	7	68	NBS75K.1	3599	69(L)
Naphthale	ene			CAS	S #: 91-20-3		
32.110	3768250	30	30	97	NBS75K.1	65151	69

#### QC Flag Legend

L - Operator selected an alternate library search match.

333-MW02-1

Lab Name: CH2M HILL

Contract: MB370

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Lab Code: MGM Case No.: MB370 SAS No.:

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SDG No.: MB370

Matrix: (soil/water) WATER

ER Lab Sample ID: MB370002

Sample wt/vol:

5.0 (g/mL) ML

Lab File ID: 18JUL1001010.D

Level: (low/med) LOW Date Received: 07/13/96

% Moisture: not dec.

Date Analyzed: 07/18/96

Column: (pack/cap) CAP

Dilution Factor: 1.0

CONCENTRATION UNITS:

CAS NO. COMPOUND (ug/L or ug/Kg) UG/L Q

·			
74-87-3	Chloromethane	10	Ū
	Vinyl chloride	10	Ū
74-83-9	Bromomethane	10	U
75-00-3	Chloroethane	10	U
75-69-4	Trichlorofluoromethane	10	ប
75-35-4 <u>-</u>	1,1-Dichloroethene	10	ט
75-09-2	Methylene chloride	10	U
156-60-5	trans-1,2-Dichloroethene	10	U
	1,1-Dichloroethane	10	U
	Chloroform	10	ט
71-55-6	1,1,1-Trichloroethane	10	Ū
56-23-5 <b>-</b>	Carbon tetrachloride	10	U
71-43-2	Benzene	10	U
	1,2-Dichloroethane	10	Ū
	Trichloroethene	10	U
78-87-5	1,2-Dichloropropane	10	U
75-27-4- <del></del>	Bromodichloromethane	10	U
110-75-8	2-Chloroethylvinyl ether	10	Ū
10061-01-5	cis-1,3-Dichloropropene	10	ט
108-88-3		10	ប
10061-02-6	trans-1,3-Dichloropropene	10	U
79-00-5	1,1,2-Trichloroethane	10.	ט
	Tetrachloroethene	10	ט
124-48-1	Dibromochloromethane	10	ט
	Chlorobenzene	10	ט
	Ethylbenzene	10	ט
	Bromoform_	10	ט
79-34-5	1,1,2,2-Tetrachloroethane	10	Ŭ

#### 1E VOLATILE ORGANICS ANALYSIS DATA SHEET TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

333-MW02-1

Lab Name: CH2M HILL

Contract: MB370

Lab Code: MGM

Case No.: MB370

LOW

SAS No.:

· SDG No.: MB370

Matrix: (soil/water) WATER

Lab Sample ID: MB370002

Sample wt/vol: 5.0

(g/mL) ML

Lab File ID: 18JUL1001010.D

Level:

(low/med)

Date Received: 07/13/96

% Moisture: not dec.

Number TICs found: 0

Data Analyzed: 07/18/96

Column: (pack/cap) CAP

Dilution Factor: 1.0

CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	0
	=======================================			
2				
4.				
5.				
/ •				
9.				·
11:				
12.	·			
14				
16.				
17.				
20.		_		
21.				
43.				
24.		_		
27.				
28.				
30.				

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Report Date: 29-Jul-1996 15:29

#### QAL, Inc:

VOLATILE REPORT METHOD 8260/5-ML PURGE

Data file : /chem/ms5.i/a071896a.b/18Jul1001010.d/

Client Smp ID: 333-MW02-1 / Autotune Date: 18-Jul-96 11:10 Lab Smp Id: MB370002 Inj Date : 18-JUL-1996 20:12 Operator : WLH/RLW

Inst ID: ms5.i

Smp Info : 333-MW02-1 MB370002

Misc Info : 5MLS /

Comment

: /chem/ms5.i/a071896a.b/8260w5.m Method

Meth Date : 29-Jul-1996 15:28 ms5 Quant Type: ISTD

Cal File: 10Jul2001020.d

Cal Date : 10-JUL-1996 22:29
Als bottle: 10
Dil Factor: 1.000
Integrator: HP RTE Compound Sublist: BROWN&ROOT.sub /

Target Version: 3.10

			CONCENTRATIONS	
	-	QUANT SIG	ON-COLUMN FINAL	L
C	ompounds	MASS	RT EXP RT REL RT RESPONSE ( ug/L) ( ug/l	L)
*	***************************************	****	秦耳 通知日区至7 四字军马马克 建合金油基苯戊苯 建含金属医苯茚 军马万二年	<b>=</b> =
•	1 Pentafluorobenzene	168.00	13.752 13.756 (1.000) 1699238 50	
	3 Chloromethane	50.00	Compound Not Detected.	
	4 Vinyl chloride	62.00	Compound Not Detected.	
	5 Bromomethane	94.00	Compound Not Detected.	
	6 Chloroethane	64.00	Compound Not Detected.	
	7 Trichlorofluoromethane	101.00	Compound Not Detected.	
	11 1,1-Dichloroethene	96.00	Compound Not Detected.	
	16 Methylene chloride	84.00	Compound Not Detected.	
	19 trans-1,2-Dichloroethene	96.00	Compound Not Detected.	
	21 1,1-Dichloroethane	63.00	Compound Not Detected.	
	31 Chloroform	83.00	Compound Not Detected.	
	33 1,1,1-Trichloroethane	97.00	Compound Not Detected.	
	35 Carbon tetrachloride	117.00	Compound Not Detected.	
*	36 1,4-Difluorobenzene	114.00	16.044 16.038 (1.000) 2188187 50	
	37 1,2-Dichloroethane	62.00	Compound Not Detected.	
	38 Benzene	78.00	Compound Not Detected.	
	39 Trichloroethene	95.00	Compound Not Detected.	
	41 1,2-Dichloropropane	63.00	Compound Not Detected.	
	43 Bromodichloromethane	83.00	Compound Not Detected.	
	45 2-Chloroethylvinyl ether	63.00	Compound Not Detected.	
	46 cis-1,3-Dichloropropene	75.00	Compound Not Detected.	
*	48 Chlorobenzene-d5	117.00	22.388 22.372 (1.000) 1706121 50	
`	49 Toluene	91.00	Compound Not Detected.	
	51 trans-1,3-Dichloropropene	75.00	Compound Not Detected.	
	52 1,1,2-Trichloroethane	83.00	Compound Not Detected.	
	55 Tetrachloroethene	166.00	Compound Not Detected.	
	56 Dibromochloromethane	129.00	Compound Not Detected.	
	58 Chlorobenzene	112.00	Compound Not Detected.	
	60 Ethylbenzene	91.00	Compound Not Detected.	

Pur mix 4/5/94
-1/29/96

Data File: /chem/ms5.i/a071896a.b/18Jul1001010.d Report Date: 29-Jul-1996 15:29

Page 2

						CONCENTRA	ATIONS
		QUANT SIG				on-Column	FINAL
C	ompounds	MASS	RT	EXP RT REL RT	RESPONSE	( ug/L)	( ug/L)
=	*************	***					
	66 Bromoform	173.00	Con	pound Not Detect	ed.		
	68 1,1,2,2-Tetrachloroethane	83.00	Con	pound Not Detect	ed.		
*	69 1,4-Dichlorobenzene-d4	152.00	27.326	27.320 (1.000)	863991	50	
\$	91 Dibromofluoromethane	113.00	13.762	13.766 (1.001)	918645	52	52
\$	92 1,2-Dichloroethane-d4	65.00	15.021	15.005 (0.936)	377640	45	45
\$	93 Toluene-d8	98.00	19.290	19.274 (1.202)	1957658	48	48
\$	94 Bromofluorobenzene	174.00	24.896	24.890 (1.112)	808644	48	48

Data File: /chem/ms5.i/a071896a.b/18Jul1001010.d

Report Date: 29-Jul-1996 16:10

Page 1

## Quality Analytical Laboratories

Unknown Compounds Quantitation Report

Data file : /chem/ms5.i/a071896a.b/18Jul1001010.d

Lab Smp Id: MB370002 Client Smp ID: 333-MW02-1

Autotune Date: 18-Jul-96 11:10:3 Inj Date : 18-JUL-1996 20:12

Operator : WLH/RLW Smp Info : 333-MW02-1 MB370002 Inst ID: ms5.i

Misc Info : 5MLS

Comment

: /chem/ms5.i/a071896a.b/8260w5.m Method

Meth Date : 29-Jul-1996 16:09 whall

Cal Date : 10-JUL-1996 22:29 Cal File: 10Jul2001020.d

Als bottle: 10

Target Version: 3.12

Dil Factor: 1.000 Integrator: HP RTE Sample Matrix: WATER Compound Sublist: BROWN&ROOT.sub

Quantitative Mode : Use RF of Nearest Std

- NO TENTATIVELY IDENTIFIED COMPOUNDS -

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333-MW03-1

Lab Name: CH2M HILL

Contract: MB370

Lab Code: MGM

Case No.: MB370

SAS No.:

SDG No.: MB370

Matrix: (soil/water) WATER

Lab Sample ID: MB370003

Sample wt/vol:

5.0 (q/mL) ML

Lab File ID:

19JUL0401004.D

Level: (low/med)

% Moisture: not dec.

LOW

Date Received: 07/13/96

Date Analyzed: 07/19/96

Column: (pack/cap) CAP

Dilution Factor: 1.0

CONCENTRATION UNITS:

CAS NO. COMPOUND (ug/L or ug/Kg) UG/L Q

·		
74-87-3Chloromethane	10	U
75-01-4Vinyl chloride	10	บ
74-83-9Bromomethane	10	انّ
75-00-3Chloroethane	10	ָ טֿ
75-69-4Trichlorofluoromethane	10	บั
75-35-41,1-Dichloroethene	10	Ŭ
75-09-2Methylene chloride	10	ן ט
156-60-5trans-1,2-Dichloroethene	10	υ
75-34-31,1-Dichloroethane	10	บ็
67-66-3Chloroform	10	U U
71-55-61,1,1-Trichloroethane	10	นี
56-23-5Carbon tetrachloride	10	บ
71-43-2Benzene		ט
	10	_
107-06-21,2-Dichloroethane	10	U
	10	Ŭ
78-87-51,2-Dichloropropane	10	Ū
75-27-4Bromodichloromethane	10	<u>ת</u>
110-75-82-Chloroethylvinyl ether	10	U
10061-01-5cis-1,3-Dichloropropene	10	Ŭ
108-88-3Toluene	10	ָּט
10061-02-6trans-1,3-Dichloropropene	10	U
79-00-51,1,2-Trichloroethane	10,	Ū
127-18-4Tetrachloroethene	10	U
124-48-1Dibromochloromethane	10	ן ט
108-90-7Chlorobenzene	10	U
100-41-4Ethylbenzene	10	ָ ט
75-25-2Bromoform	10	บ
79-34-51,1,2,2-Tetrachloroethane	10	U
	<u> </u>	

# VOLATILE ORGANICS ANALYSIS DATA SHEET TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

333-MW03-1

Lab Name: CH2M HILL

Contract: MB370

Lab Code: MGM Case No.: MB370

SAS No.:

SDG No.: MB370

Matrix: (soil/water) WATER

Lab Sample ID: MB370003

Sample wt/vol: 5.0

(g/mL) ML

Lab File ID:

19JUL0401004.D

Level: (low/med) LOW

Date Received: 07/13/96

% Moisture: not dec.

Data Analyzed: 07/19/96

Column: (pack/cap) CAP

Number TICs found: 0

Dilution Factor: 1.0

CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.				
3				
5				
6. 7.	· · · · · · · · · · · · · · · · · · ·	_		
9.				
11:				
12.		_		
14.				
16		_		
19		_		
21.		_		
23.				
25.				
28.	***************************************			
29.				

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SW846-8260

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Report Date: 29-Jul-1996 16:02

#### Quality Analytical Laboratories

VOLATILE REPORT METHOD 8260/5-ML PURGE

Data file : /chem/ms5,i/a071996a.b/19Jul0401004.d/

Client Smp ID: 333-MW03-1 Lab Smp Id: MB370003 / Autotune Date: 19-Jul-96 09:48

Inj Date : 19-JUL-1996 11:57 Operator : WLH/RLW Smp Info : 333-MW03-1 MB370003 Inst ID: ms5.i ✓

Misc Info : 5MLS /

Comment

: /chem/ms5.i/a071996a.b/8260w5.m / Method

Meth Date : 29-Jul-1996 16:02 whall Quant Type: ISTD

Cal Date : 10-JUL-1996 22:29 Cal File: 10Jul2001020.d /

Als bottle: 4
Dil Factor: 1.000 /
Integrator: HP RTE

Compound Sublist: BROWN&ROOT.sub /

Target Version: 3.12

Concentration Formula: Uf \* 5/Vo

Name	Value	Description				
Uf	1.000	ng unit correction factor				
Vo	5.000	Sample Volume purged (mL)				

								CONCENT	RATIONS
			QUANT SIG					ON-COLUMN	FINAL
Compounds		ınds	MASS	RT	EXP RT	REL RT RESP	ONSE	( ug/L)	( ug/L)
看着进足双军等往 油鱼采菜 耳朵 化苯甲基苯基苯基		, , , , , , , , , , , , , , , , , , ,	<del>-</del>	==	e=====			*****	*****
*	1	Pentafluorobenzene	168.00	13.764	13.724	(1.000) 169	4298	50	
•	3	Chloromethane	50.00		3.199	Compound	Not I	Detected.	
	4	Vinyl chloride	62.00		3.524	Compound	Not I	Detected.	
	5	Bromomethane	94.00		4.733	Compound	Not I	Detected.	
	6	Chloroethane	64.00		5.048	Compound	Not I	Detected.	
	7	Trichlorofluoromethane	101.00		5.806	Compound	Not I	Detected.	
	11	1,1-Dichloroethene	96.00		7.566	Compound	Not I	Detected.	
	16	Methylene chloride	84.00		9.012	Compound	Not I	Detected.	
	19	trans-1,2-Dichloroethene	96.00		9.809	Compound	Not I	Detected.	
	21	1,1-Dichloroethane	63.00		11.058	Compound	Not I	Detected.	
	31	Chloroform	83.00		13.163	Compound	Not I	Detected.	
	33	1,1,1-Trichloroethane	97.00		14.206	Compound	Not I	Detected.	
	35	Carbon tetrachloride	117.00		14.786	Compound	Not I	Detected.	
*	36	1,4-Difluorobenzene	114.00	16.046	16.025	(1.000) 216	0075	50	*
	37	1,2-Dichloroethane	62.00		15.219	Compound	Not I	Detected.	
	38	Benzene	78.00		15.189	Compound	Not I	Detected.	
	39	Trichloroethene	95.00		16.665	Compound	Not I	Detected.	
	41	1,2-Dichloropropane	63.00		17.058	Compound	Not I	Detected.	
	43	Bromodichloromethane	83.00		17.560	Compound	Not I	Detected.	
	45	2-Chloroethylvinyl ether	63.00		18.386	Compound	Not I	Detected.	-
	46	cis-1,3-Dichloropropene	75.00		18.799	Compound	Not I	Detected.	
•	48	Chlorobenzene-d5	117.00	22.381	22.370	(1.000) 166	9608	50	
	49	Toluene	91.00		19.439	Compound	Not I	Detected.	

mp 8/2/96

								CONCENTRA	ATIONS
			QUANT SIG					ON-COLUMN	FINAL
Compounds		unds	MASS	RT	EXP RT	REL RT	RESPONSE	( ug/L)	( ug/L)
					****				******
	51	trans-1,3-Dichloropropene	75.00		19.871	Comp	pound Not De	etected.	
	52	1,1,2-Trichloroethane	83.00		20.157	Comp	pound Not De	etected.	
	55	Tetrachloroethene	166.00		20.776	Comp	pound Not De	etected.	
	56	Dibromochloromethane	129.00		21.160	Com	pound Not De	tected.	
	58	Chlorobenzene	112.00		22.458	Comp	pound Not De	etected.	
	60	Ethylbenzene	91.00		22.606	Comp	pound Not De	etected.	
	66	Bromoform	173.00		24.337	Comp	pound Not De	etected.	
	68	1,1,2,2-Tetrachloroethane	83.00		24.760	Comp	pound Not De	etected.	
*	69	1,4-Dichlorobenzene-d4	152.00	27.328	27.327	(1.000)	829303	50	•
\$	91	Dibromofluoromethane	113.00	13.764	13.743	(1.000)	861076	48	48
\$	92	1,2-Dichloroethane-d4	65.00	15.013	14.993	(0.936)	329439	40	40
\$	93	Toluene-d8	98.00	19.282	19.271	(1.202)	1878255	47	47
s	94	Bromofluorobenzene	174.00	24.899	24.888	(1.113)	751757	46	46

Data File: /chem/ms5.i/a071996a.b/19Jul0401004.d

Report Date: 29-Jul-1996 16:02

## Quality Analytical Laboratories

Unknown Compounds Quantitation Report

Data file : /chem/ms5.i/a071996a.b/19Jul0401004.d

Lab Smp Id: MB370003 Client Smp ID: 333-MW03-1

Inj Date : 19-JUL-1996 11:57 Autotune Date: 19-Jul-96 09:48:4

Operator : WLH/RLW Inst ID: ms5.i

Smp Info : 333-MW03-1 MB370003 Misc Info : 5MLS

Comment

Method : /chem/ms5.i/a071996a.b/8260w5.m Meth Date : 29-Jul-1996 16:02 whall

Cal Date : 10-JUL-1996 22:29 Cal File: 10Jul2001020.d

Als bottle: 4

Dil Factor: 1.000 Target Version: 3.12

Integrator: HP RTE Compound Sublist: BROWN&ROOT.sub

Sample Matrix: WATER

Quantitative Mode : Use RF of Nearest Std

- NO TENTATIVELY IDENTIFIED COMPOUNDS -

Page 3

# VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO

333-PC4-1

Lab Name: CH2M HILL

Contract: MB370

Lab Code: MGM

Case No.: MB370

SAS No.:

SDG No.: MB370

Matrix: (soil/water) WATER

Lab Sample ID: MB370004

Sample wt/vol:

5.0 (g/mL) ML

Lab File ID:

19JUL0701007.D

Level: (low/med)

Date Received: 07/13/96

% Moisture: not dec.

Date Analyzed: 07/19/96

Column: (pack/cap) CAP

Dilution Factor: 1.0

CONCENTRATION UNITS:

CAS NO. COMPOUND (ug/L or ug/Kg) UG/L

LOW

Q

		<del></del> ,
74-87-3Chloromethane	10	บ
75-01-4Vinyl chloride	10	Ŭ
74-83-9Bromomethane	10	บ
75-00-3Chloroethane	10	Ü
75-69-4Trichlorofluoromethane	10	ָ <u>ט</u>
75-35-41,1-Dichloroethene		Ü
75-09-2Methylene chloride		บ
156-60-5trans-1,2-Dichloroethene		บ
75-34-31,1-Dichloroethane	10	Ü
67-66-3Chloroform	10	الق
71-55-61,1,1-Trichloroethane	10	ש
56-23-5Carbon tetrachloride	10	Ü
71-43-2Benzene	10	-
	10	Ŭ
107-06-21,2-Dichloroethane 79-01-6Trichloroethene	. [	Ŭ
	10	Ü
78-87-51,2-Dichloropropane	. 10	ן דַ
75-27-4Bromodichloromethane	. 10	Ü
110-75-82-Chloroethylvinyl ether	. 10	ָ טַ
10061-01-5cis-1,3-Dichloropropene	. 10	<u>ט</u>
108-88-3Toluene	. 10	Ū
10061-02-6trans-1,3-Dichloropropene	. 10	ע
79-00-51,1,2-Trichloroethane	. 10.	<u>ט</u>
127-18-4Tetrachloroethene	.  10	ן ט
124-48-1Dibromochloromethane	.  10	ן ט
108-90-7Chlorobenzene	10	ָ <u>ע</u>
100-41-4Ethylbenzene	. 10	ע
75-25-2Bromoform	.  10	ן ט
79-34-51,1,2,2-Tetrachloroethane	10	Ū
	.	

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VOLATILE ORGANICS ANALYSIS DATA SHEET TENTATIVELY IDENTIFIED COMPOUNDS

333-PC4-1

Lab Name: CH2M HILL

Contract: MB370

EPA SAMPLE NO

Lab Code: MGM Case No.: MB370

SAS No.: SDG No.: MB370

Matrix: (soil/water) WATER

Lab Sample ID: MB370004

Sample wt/vol: 5.0 (g/mL) ML

Lab File ID: 19JUL0701007.D

Level: (low/med) LOW

Date Received: 07/13/96

% Moisture: not dec.

Data Analyzed: 07/19/96

Column: (pack/cap) CAP

Dilution Factor: 1.0

CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L

Number TICs found: 0

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.				
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J		-	<del>-                                   </del>	
5	<del></del>	-		
6.				
/ ·		-		
8 . 9 .	<del></del>	-	<del></del>	
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12.		-		
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15. I		.		
16:		-		
18.				
19.				
20.	······································	-		
22.		-		
23.				
24.				
25. 26.	<del></del>	-		
27.				
28.				
29.		-		

SW846-8260

Report Date: 29-Jul-1996 16:03

Page 1

### Quality Analytical Laboratories

VOLATILE REPORT METHOD 8260/5-ML PURGE

Data file : /chem/ms5.i/a071996a.b/19Jul0701007.d/

Lab Smp Id: MB370004 / Client Smp ID: 333-PC4-1 —

Inj Date : 19-JUL-1996 14:02 Operator : WLH/RLW Autotune Date: 19-Jul-96 09:48

Inst ID: ms5.i

Smp Info : 333-PC4-1 MB370004

Misc Info : 5MLS /

Comment

: /chem/ms5.i/a071996a.b/8260w5.m/ Method

Meth Date : 29-Jul-1996 16:02 whall Quant Type: ISTD

Cal Date: 10-JUL-1996 22:29
Als bottle: 7
Dil Factor: 1.000 Cal File: 10Jul2001020.d /

Integrator: HP RTE Compound Sublist: BROWN&ROOT.sub \_

Target Version: 3.12

Concentration Formula: Uf \* 5/Vo

Name	Value	Description
U£ .	1.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)

								CONCENTR	ATIONS
			QUANT SIG					ON-COLUMN	FINAL
,n	ιροι	ınds	MASS	RT	EXP RT	REL RT RES	PONSE	( ug/L)	( ug/L)
===	* * * *	, s 3 x 2 2 2 2 2 2 2 2 2 2 2 2 2 2 2 2 2 2	====	==	22222	E33333 =33		======	2333223
*	1	Pentafluorobenzene	168.00	13.761	13.724	(1.000) 16	48174	50	
	3	Chloromethane	50.00	•	3.199	Compound	Not	Detected.	
	4	Vinyl chloride	62.00		3.524	Compound	Not	Detected.	
	5	Bromomethane	94.00		4.733	Compound	Not	Detected.	
	6	Chloroethane	64.00		5.048	Compound	Not	Detected.	
	7	Trichlorofluoromethane	101.00		5.806	Compound	Not	Detected.	
	11	1,1-Dichloroethene	96.00		7.566	Compound	Not	Detected.	
	16	Methylene chloride	84.00		9.012	Compound	Not	Detected.	
	19	trans-1,2-Dichloroethene	96.00		9.809	Compound	Not	Detected.	
	21	1,1-Dichloroethane	63.00		11.058	Compound	Not	Detected.	
	31	Chloroform	83.00		13.163	Compound	Not	Detected.	
	33	1,1,1-Trichloroethane	97.00		14.206	Compound	Not	Detected.	
	35	Carbon tetrachloride	117.00		14.786	Compound	Not	Detected.	
*	36	1,4-Difluorobenzene	114.00	16.043	16.025	(1.000) 21	30537	50	
	37	1,2-Dichloroethane	62.00		15.219	Compound	Not	Detected.	
	38	Benzene	78.00		15.189	Compound	Not	Detected.	
	39	Trichloroethene	95.00		16.665	Compound	Not	Detected.	
	41	1,2-Dichloropropane	63.00		17.058	Compound	Not	Detected.	
	43	Bromodichloromethane	83.00		17.560	Compound	Not	Detected.	
	45	2-Chloroethylvinyl ether	63.00		18.386	Compound	Not	Detected.	
	46	cis-1,3-Dichloropropene	75.00		18.799	Compound	Not	Detected.	
*	48	Chlorobenzene-d5	117.00	22.397	22.370	(1.000) 16	55637	50	
	49	Toluene	91.00		19.439	Compound	Not	Detected.	

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							CONCENTRA	TIONS
	•	QUANT SIG					ON-COLUMN	FINAL
Co	ompounds	MASS	RT	EXP RT	REL RŤ	RESPONSE	( ug/L)	( ug/L)
-					*****			******
	51 trans-1,3-Dichloropropene	75.00	•	19.871	Comp	ound Not De	etected.	
	52 1,1,2-Trichloroethane	83.00		20.157	Comp	ound Not De	etected.	
	55 Tetrachloroethene	166.00		20.776	Comp	ound Not De	etected.	
	56 Dibromochloromethane	129.00		21.160	Comp	ound Not De	tected.	
	58 Chlorobenzene	112.00		22.458	Comp	ound Not De	tected.	
	60 Ethylbenzene	91.00		22.606	Comp	ound Not De	tected.	
	66 Bromoform	173.00		24.337	Comp	ound Not De	tected.	
	68 1,1,2,2-Tetrachloroethane	83.00		24.760	Comp	ound Not De	tected.	
*	69 1,4-Dichlorobenzene-d4	152.00	27.335	27.327	(1.000)	828235	50	
\$	91 Dibromofluoromethane	113.00	13.771	13.743	(1.001)	921140	53	53
\$	92 1,2-Dichloroethane-d4	65.00	15.020	14.993	(0.936)	369928	45	45
\$	93 Toluene-d8	98.00	19.289	19.271	(1.202)	1957457	49	49
s	94 Bromofluorobenzene	174.00	24.906	24.888	(1.112)	804497	49	49

Data File: /chem/ms5.i/a071996a.b/19Jul0701007.d Page 3

Report Date: 29-Jul-1996 16:03

### Quality Analytical Laboratories

Unknown Compounds Quantitation Report

Data file : /chem/ms5.i/a071996a.b/19Jul0701007.d

Lab Smp Id: MB370004 Client Smp ID: 333-PC4-1

Inj Date : 19-JUL-1996 14:02 Autotune Date: 19-Jul-96 09:48:4

Operator : WLH/RLW Inst ID: ms5.i

Smp Info : 333-PC4-1 MB370004

Misc Info : 5MLS

Comment

: /chem/ms5.i/a071996a.b/8260w5.m Method

Meth Date : 29-Jul-1996 16:02 whall

Cal Date : 10-JUL-1996 22:29 Cal File: 10Jul2001020.d

Als bottle: 7

Dil Factor: 1.000 Integrator: HP RTE Target Version: 3.12

Compound Sublist: BROWN&ROOT.sub

Sample Matrix: WATER

Quantitative Mode : Use RF of Nearest Std

- NO TENTATIVELY IDENTIFIED COMPOUNDS -

5 . 1

# VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

333-MW01-1B

Lab Name: CH2M HILL

Contract: MB370

SAS No.:

Lab Code: MGM Case No.: MB370

LOW

SDG No.: MB370

Matrix: (soil/water) WATER

Lab Sample ID: MB370005

Sample wt/vol:

5.0 (g/mL) ML

Lab File ID: 19JUL0501005.D

Level: (low/med)

Date Received: 07/13/96

% Moisture: not dec.

Date Analyzed: 07/19/96

Column: (pack/cap) CAP

Dilution Factor: 1.0

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/L CAS NO. COMPOUND Q

74-87-3	Chloromethane	10	บ
	Vinyl chloride	10	ן מ
	Bromomethane	10	ש
	Chloroethane	10	ן ט
	Trichlorofluoromethane	10	ט
	1,1-Dichloroethene	10	ט
	Methylene chloride	10	υ
	trans-1,2-Dichloroethene	10	ט
	1,1-Dichloroethane	10	ן ט
	Chloroform	10	ប
7155-6	1,1,1-Trichloroethane	10	υ
	Carbon tetrachloride	10	ט
71-43-2		10	ט
107-06-2	1, 2-Dichloroethane	10	ט
79-01-6	Trichloroethene	10	ַ ט
78-87-5	1,2-Dichloropropane	10	Ū
	Bromodichloromethane	10	ט
110-75-8	2-Chloroethylvinyl ether	10	U
10061-01-5	cis-1,3-Dichloropropene	10	U
108-88-3		10	U
	trans-1,3-Dichloropropene	10	ן ט
79-00-5	1,1,2-Trichloroethane	10-	ַ ט
	Tetrachloroethene	10	ן ע
124-48-1	Dibromochloromethane	10	ן ט
	Chlorobenzene	. 10	ָ ע
100-41-4	Ethylbenzene	10	U
75-25-2	Bromoform	10	U
79-34-5	1,1,2,2-Tetrachloroethane	10	ט (
		l	l

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VOLATILE ORGANICS ANALYSIS DATA SHEET TENTATIVELY IDENTIFIED COMPOUNDS

333-MW01-1B

19JUL0501005.D

Lab File ID:

Data Analyzed: 07/19/96

EPA SAMPLE NO.

Lab Name: CH2M HILL Contract: MB370

Sample wt/vol: 5.0

% Moisture: not dec.

SDG No.: MB370 Lab Code: MGM Case No.: MB370 SAS No.:

Matrix: (soil/water) WATER Lab Sample ID: MB370005

Level: (low/med) LOW Date Received: 07/13/96

Column: (pack/cap) CAP Dilution Factor: 1.0

CONCENTRATION UNITS: Number TICs found: 0 (ug/L or ug/Kg) UG/L

(q/mL) ML

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q =====
1				
٠				
5.				
7				
9:				
10.				
13.				
15.				
16:				
18.				
21.				
22.				
24.				
26.				
28. 29. 30.				
30				

SW846-8260

N . 4

Report Date: 29-Jul-1996 16:02

### Quality Analytical Laboratories

VOLATILE REPORT METHOD 8260/5-ML PURGE

Data file : /chem/ms5.i/a071996a.b/19Jul0501005.d/

Lab Smp Id: MB370005 -Client Smp ID: 333-MW01-1B -Inj Date : 19-JUL-1996 12:39 Autotune Date: 19-Jul-96 09:48

Operator : WLH/RLW Inst ID: ms5.i

Smp Info : 333-MW01-1B MB370005

Misc Info : 5MLS /

Comment

: /chem/ms5.i/a071996a.b/8260w5.m / Method

Meth Date : 29-Jul-1996 16:02 whall Cal Date : 10-JUL-1996 22:29 Quant Type: ISTD

Cal File: 10Jul2001020.d

Als bottle: 5

Dil Factor: 1.000 -

Integrator: HP RTE Compound Sublist: BROWN&ROOT.sub ~

Target Version: 3.12

Concentration Formula: Uf \* 5/Vo

Name	Value	Description
Uf	1.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)

								CONCENTRA	ATIONS
			QUANT SIG					ON-COLUMN	FINAL
nc.	poi	ınds	MASS	RT	EXP RT	REL RT RES	PONSE	( ug/L)	( ug/L)
		****************	***	= =	*****	****** 7==		33332E	======
*	1	Pentafluorobenzene	168.00	13.764	13.724	(1.000) 16	45948	50	
	3	Chloromethane	50.00		3.199	Compound	Not	Detected.	
	4	Vinyl chloride	62.00		3.524	Compound	Not	Detected.	
	5	Bromomethane	94.00		4.733	Compound	Not	Detected.	
	6	Chloroethane	64.00		5.048	Compound	Not	Detected.	
	7	Trichlorofluoromethane	101.00		5.806	Compound	Not	Detected.	
	11	1,1-Dichloroethene	96.00		7.566	Compound	Not	Detected.	
	16	Methylene chloride	84.00		9.012	Compound	Not	Detected.	
	19	trans-1,2-Dichloroethene	96.00		9.809	Compound	Not	Detected.	
	21	1,1-Dichloroethane	63.00		11.058	Compound	Not	Detected.	
	31	Chloroform	83.00		13.163	Compound	Not	Detected.	
	33	1,1,1-Trichloroethane	97.00		14.206	Compound	Not	Detected.	
	35	Carbon tetrachloride	117.00		14.786	Compound	Not	Detected.	
*	36	1,4-Difluorobenzene	114.00	16.046	16.025	(1.000) 21	.22937	50	
	37	1,2-Dichloroethane	62.00		15.219	Compound	Not	Detected.	
	38	Benzene	78.00		15.189	Compound	Not	Detected.	
	39	Trichloroethene	95.00		16.665	Compound	Not	Detected.	
	41	1,2-Dichloropropane	63.00		17.058	Compound	Not	Detected.	
	43	Bromodichloromethane	83.00		17.560	Compound	Not	Detected.	
	45	2-Chloroethylvinyl ether	63.00		18.386	Compound	Not	Detected.	
	46	cis-1,3-Dichloropropene	75.00		18.799	Compound	Not	Detected.	
٠	48	Chlorobenzene-d5	117.00	22.391	22.370	(1.000) 16	41119	50	

Data File: /chem/ms5.i/a071996a.b/19Jul0501005.d Report Date: 29-Jul-1996 16:02

							CONCENTRA	ATIONS	
		QUANT SIG					on-column	FINAL	
Compo	ounds	MASS	RT	EXP RT	REL RT	RESPONSE	( ug/L)	( ug/L)	
					*****		******		below PQL
4 9	9 Toluene	91.00	19.450	19.439	(0.869)	60197	1	الملار	METOLOGIA
5	1 trans-1,3-Dichloropropene	75.00		19.871	Com	pound Not De	tected.		ماهار
5	2 1,1,2-Trichloroethane	83.00		20.157	Com	pound Not De	etected.		1/3/1/10
5	5 Tetrachloroethene	166.00		20.776	Com	pound Not De	etected.		·
56	5 Dibromochloromethane	129.00		21.160	Com	pound Not De	etected.		
51	8 Chlorobenzene	112.00		22.458	Com	pound Not De	etected.		
6	O Ethylbenzene	91.00		22.606	Соп	pound Not De	etected.		
6	6 Bromoform	173.00		24.337	Com	pound Not De	etected.		
68	3 1,1,2,2-Tetrachloroethane	83.00		24.760	Соп	pound Not De	etected.		
* 69	9 1,4-Dichlorobenzene-d4	152.00	27.328	27.327	(1.000)	824309	50		
\$ 9:	1 Dibromofluoromethane	113.00	13.764	13.743	(1.000)	892257	52	52	
\$ 93	2 1,2-Dichloroethane-d4	65.00	15.013	14.993	(0.936)	363691	44	44	
\$ 93	3 Toluene-d8	98.00	19.292	19.271	(1.202)	1925246	49	49	
\$ 94	4 Bromofluorobenzene	174.00	24.899	24.888	(1.112)	787391	49	49	

## QC Flag Legend

a - Target compound detected but, quantitated amount Below Limit Of Quantitation(BLOQ).

Data File: /chem/ms5.i/a071996a.b/19Jul0501005.d

Report Date: 29-Jul-1996 16:02

### Quality Analytical Laboratories

Unknown Compounds Quantitation Report

Data file : /chem/ms5.i/a071996a.b/19Jul0501005.d

Lab Smp Id: MB370005 Client Smp ID: 333-MW01-1B

Autotune Date: 19-Jul-96 09:48:4

Inj Date : 19-JUL-1996 12:39 Operator : WLH/RLW Smp Info : 333-MW01-1B MB370005 Inst ID: ms5.i

Misc Info : 5MLS

Comment

: /chem/ms5.i/a071996a.b/8260w5.m Method

Meth Date : 29-Jul-1996 16:02 whall

Cal Date : 10-JUL-1996 22:29 Cal File: 10Jul2001020.d

Als bottle: 5

Target Version: 3.12

Dil Factor: 1.000 Integrator: HP RTE Compound Sublist: BROWN&ROOT.sub

Sample Matrix: WATER

Ouantitative Mode : Use RF of Nearest Std

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Page 3

<sup>-</sup> NO TENTATIVELY IDENTIFIED COMPOUNDS -

# VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO

333-MW03-1D

Lab Name: CH2M HILL

Contract: MB370

SDG No.: MB370 SAS No.:

Matrix: (soil/water) WATER

Lab Sample ID: MB370006

Sample wt/vol:

5.0 (g/mL) ML

Case No.: MB370

LOW

Lab File ID:

19JUL0601006.D

Level: (low/med)

Lab Code: MGM

Date Received: 07/13/96

% Moisture: not dec.

Date Analyzed: 07/19/96

÷ ,:

Column: (pack/cap) CAP

Dilution Factor: 1.0

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/L CAS NO. COMPOUND Q

·			
74-87-3	Chloromethane	10	υ
	Vinyl chloride	10	ט
74-83-9		10	ט
75-00-3		10	ן ט
	Trichlorofluoromethane	10	ט
75-35-4	1,1-Dichloroethene	10	ש
75-09-2	Methylene chloride	10	ט
156-60-5	trans-1,2-Dichloroethene	10	ן . טן
	1,1-Dichloroethane	10	ן ט
67-66-3		10	ប
71-55-6	1,1,1-Trichloroethane	10	ן ט
56-23-5	Carbon tetrachloride	10	ן ט
71-43-2		10	ן ט
107-06-2	1,2-Dichloroethane	10	ן די
	Trichloroethene	10	ן די
	1,2-Dichloropropane	10	ן ע
	Bromodichloromethane	10	ן ט
110-75-8	2-Chloroethylvinyl ether	10	ן ט
	cis-1,3-Dichloropropene	10	ן ט
108-88-3		10	ן ט
	trans-1,3-Dichloropropene	10	ן די ן
	1,1,2-Trichloroethane	10-	ן די ן
127-18-4	Tetrachloroethene	10	U
124-48-1	Dibromochloromethane	10	ן די
	Chlorobenzene	10	ן די
100-41-4	Ethylbenzene	10	1 3
75-25-2		10	ן ט
79-34-5	1,1,2,2-Tetrachloroethane	10	ן די

# VOLATILE ORGANICS ANALYSIS DATA SHEET TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

333-MW03-1D

Lab Name: CH2M HILL

Contract: MB370

Lab Code: MGM

Case No.: MB370

SAS No.:

SDG No.: MB370

Matrix: (soil/water) WATER

Lab Sample ID: MB370006

Lab File ID:

19JUL0601006.D

Sample wt/vol: 5.0 Level: (low/med)

Number TICs found: 0

(g/mL) ML

Date Received: 07/13/96

% Moisture: not dec.

Data Analyzed: 07/19/96

Column: (pack/cap) CAP

LOW

Dilution Factor: 1.0

CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1				
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4				
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<b>⊥⊥.</b>				
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Data File: /chem/ms5.i/a071996a.b/19Jul0601006.d

Report Date: 29-Jul-1996 16:02

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#### Quality Analytical Laboratories

VOLATILE REPORT METHOD 8260/5-ML PURGE

Data file : /chem/ms5.i/a071996a.b/19Jul0601006.d /

Lab Smp Id: MB370006 🗸 Client Smp ID: 333-MW03-1D / Inj Date : 19-JUL-1996 13:20 Autotune Date: 19-Jul-96 09:48

Operator : WLH/RLW Inst ID: ms5.i /

Smp Info : 333-MW03-1D MB370006

Misc Info : 5MLS/

Comment

Method : /chem/ms5.i/a071996a.b/8260w5.m /
Meth Date : 29-Jul-1996 16:02 whall Quant Type: ISTD

Cal Date : 10-JUL-1996 22:29 Cal File: 10Jul2001020.d /

Als bottle: 6

Dil Factor: 1.000/

Integrator: HP RTE Compound Sublist: BROWN&ROOT.sub /

Target Version: 3.12

Concentration Formula: Uf \* 5/Vo

Name	Vælue	Description
U£		ng unit correction factor
Vo	5.000	Sample Volume purged (mL)

						1		CONCENTRA	ATIONS
			QUANT SIG					ON-COLUMN	FINAL
Cor	npou	ınds	MASS	RT	EXP RT	REL RT RES	PONSE	( ug/L)	( ug/L)
==:		· 医罗尔里耳耳耳耳耳耳耳耳耳耳耳耳耳耳耳耳耳耳	****	*=		****** **=		****	******
1.	1	Pentafluorobenzene	168.00	13.765	13.724	(1.000) 16	53073	50	
	3	Chloromethane	50.00		3.199	Compound	Not	Detected.	
	4	Vinyl chloride	62.00		3.524	Compound	Not	Detected.	
	5	Bromomethane	94.00		4.733	Compound	Not	Detected.	
	6	Chloroethane	64.00		5.048	Compound	Not	Detected.	
	7	Trichlorofluoromethane	101.00		5.806	Compound	Not	Detected.	
	11	1,1-Dichloroethene	96.00		7.566	Compound	Not	Detected.	
	16	Methylene chloride	84.00		9.012	Compound	Not	Detected.	
	19	trans-1,2-Dichloroethene	96.00		9.809	Compound	Not	Detected.	
	21	1,1-Dichloroethane	63.00		11.058	Compound	Not	Detected.	
	31	Chloroform	83.00		13.163	Compound	Not	Detected.	
	33	1,1,1-Trichloroethane	97.00		14.206	Compound	Not	Detected.	
	35	Carbon tetrachloride	117.00		14.786	Compound	Not	Detected.	
~ <b>*</b>	36	1,4-Difluorobenzene	114.00	16.047	16.025	(1,000) 21	34849	50	
	37	1,2-Dichloroethane	62.00		15.219	Compound	Not	Detected.	
	38	Benzene	78.00		15.189	Compound	Not	Detected.	
	39	Trichloroethene	95.00		16.665	Compound	Not	Detected.	
	41	1,2-Dichloropropane	63.00		17.058	Compound	Not	Detected.	
	43	Bromodichloromethane	83.00		17.560	Compound	Not	Detected.	
	45	2-Chloroethylvinyl ether	63.00		18.386	Compound	Not	Detected.	
	46	cis-1,3-Dichloropropene	75.00		18.799	Compound	Not	Detected.	
*	48	Chlorobenzene-d5	117.00	22.391	22.370	(1.000) 16	54485	50	
	49	Toluene	91.00		19.439	Compound	Not	Detected.	

							CONCENTRA	TIONS
		QUANT SIG					ON-COLUMN	FINAL
C	ompounds	MASS	RT	EXP RT	REL RT	RESPONSE	( ug/L)	( ug/L)
-	医电对电流性 法法律证 医甲状腺 医皮肤 医皮肤 医牙术 法证明		==			******	******	****
	51 trans-1,3-Dichloropropene	75.00		19.871	Comp	ound Not De	etected.	
	52 1,1,2-Trichloroethane	83.00		20.157	Comp	ound Not De	etected.	
	55 Tetrachloroethene	166.00		20.776	Comp	ound Not De	etected.	
	56 Dibromochloromethane	129.00		21.160	Comp	ound Not De	etected.	
	58 Chlorobenzene	112.00		22.458	Comp	ound Not De	etected.	
	60 Ethylbenzene	91.00		22.606	Comp	ound Not De	etected.	
	66 Bromoform	173.00		24.337	Comp	ound Not De	etected.	
	68 1,1,2,2-Tetrachloroethane	83.00		24.760	Comp	ound Not De	etected.	
*	69 1,4-Dichlorobenzene-d4	152.00	27.329	27.327	(1.000)	829727	50	
\$	91 Dibromofluoromethane	113.00	13.775	13.743	(1.001)	912489	53	53
\$	92 1,2-Dichloroethane-d4	65.00	15.024	14.993	(0.936)	367756	45	45
\$	93 Toluene-d8	98.00	19.283	19.271	(1.202)	1934736	49	49
\$	94 Bromofluorobenzene	174.00	24.900	24.888	(1.112)	792172	48	4.8

Data File: /chem/ms5.i/a071996a.b/19Jul0601006.d Page 3

Report Date: 29-Jul-1996 16:02

### Quality Analytical Laboratories

Unknown Compounds Quantitation Report

Data file : /chem/ms5.i/a071996a.b/19Jul0601006.d

Lab Smp Id: MB370006 Client Smp ID: 333-MW03-1D

Autotune Date: 19-Jul-96 09:48:4 Inj Date : 19-JUL-1996 13:20

Inst ID: ms5.i Operator : WLH/RLW

Smp Info : 333-MW03-1D MB370006

Misc Info : 5MLS

Comment

Method : /chem/ms5.i/a071996a.b/8260w5.m Meth Date : 29-Jul-1996 16:02 whall

Cal File: 10Jul2001020.d Cal Date : 10-JUL-1996 22:29

Als bottle: 6
Dil Factor: 1.000
Integrator: HP RTE Target Version: 3.12

Compound Sublist: BROWN&ROOT.sub

Sample Matrix: WATER

Quantitative Mode : Use RF of Nearest Std

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<sup>-</sup> NO TENTATIVELY IDENTIFIED COMPOUNDS -

EPA SAMPLE NO.

TRIP\_BLANK

Q

Lab Name: CH2M HILL

Contract: MB370

Lab Code: MGM Case No.: MB370 SAS No.: SDG No.: MB370

Matrix: (soil/water) WATER Lab Sample ID: MB370007

Sample wt/vol: 5.0 (g/mL) ML Lab File ID: 18JUL0801008.D

Level: (low/med) LOW Date Received: 07/13/96

% Moisture: not dec. Date Analyzed: 07/18/96

Column: (pack/cap) CAP Dilution Factor: 1.0

CAS NO. COMPOUND CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L

74-87-3	Chloromethane	10	U
	Vinyl chloride	10	ان ا
74-83-9	Bromomethane	10	ا تا
	Chloroethane	10	ן ט
75-69-4	Trichlorofluoromethane	10	ן די
75-35-4=	1,1-Dichloroethene	10	ן די
75-09-2 <b></b>	Methylene chloride	5	J
	trans-1,2-Dichloroethene	10	ט
75-34-3	1,1-Dichloroethane	10	ן ט
67-66-3	Chloroform	10	ן ט
	1,1,1-Trichloroethane	10	ן ט
56-23-5	Carbon tetrachloride	10	ן ט
71-43-2		10	ן ט
	1,2-Dichloroethane	10	ן ט
	Trichloroethene	10	1 1
	1,2-Dichloropropane	10	U
	Bromodichloromethane	10	U
110-75-8	2-Chloroethylvinyl ether	10	ן די
10061-01-5	cis-1,3-Dichloropropene	10	ן ען
108-88-3		10	[U [
10061-02-6	trans-1,3-Dichloropropene	10	U
79-00-5	1,1,2-Trichloroethane	10,	4 1
	Tetrachloroethene	10	U
	Dibromochloromethane	10	ן די
	Chlorobenzene	10	U
	Ethylbenzene	10	ן די
	Bromoform_	10	1 1
79-34-5	1,1,2,2-Tetrachloroethane	10	ן ט
			ll

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#### 1E VOLATILE ORGANICS ANALYSIS DATA SHEET TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

TRIP BLANK

Tab Name: CH2M HILL

Contract: MB370

Lab Code: MGM Case No.: MB370 SAS No.:

SDG No.: MB370

Matrix: (soil/water) WATER

Lab Sample ID: MB370007

Sample wt/vol: 5.0 (g/mL) ML

Lab File ID: 18JUL0801008.D

Level: (low/med) LOW

Date Received: 07/13/96

% Moisture: not dec.

Number TICs found: 0

Data Analyzed: 07/18/96

Column: (pack/cap) CAP

Dilution Factor: 1.0

CONCENTRATION UNITS: (uq/L or uq/Kq) UG/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
	=======================================	======	=========	====
1. 2.				
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43.				
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20.				
29.				
		<del></del>	<del></del>	

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Report Date: 29-Jul-1996 15:28

### QAL, Inc:

VOLATILE REPORT METHOD 8260/5-ML PURGE

Data file : /chem/ms5.i/a071896a.b/18Jul0801008.d ~

Client Smp ID: TRIP BLANK -Lab Smp Id: MB370007 -Inj Date : 18-JUL-1996 18:51 Autotune Date: 18-Jul-96 11:10

Operator : WLH/RLW Inst ID: ms5.i

Smp Info : TRIP\_BLANK MB370007

Misc Info : 5MLS

Comment

Cal File: 10Jul2001020.d /

Als bottle: 8

Dil Factor: 1.000

Integrator: HP RTE Compound Sublist: BROWN&ROOT.sub

Target Version: 3.10

				CONCENTRATIONS
	-	QUANT SIG		ON-COLUMN FINAL
Co	ompounds .	MASS	RT EXP RT REL RT RESPONSE	(ug/L) $(ug/L)$
	医复数骨骨 医牙牙 医甲基甲基 地名西班牙 计多元	*===		*******
*	1 Pentafluorobenzene	168.00	13.755 13.756 (1.000) 1717586	50
	3 Chloromethane	50.00	Compound Not Detected.	
	4 Vinyl chloride	62.00	Compound Not Detected.	
	5 Bromomethane	94.00	Compound Not Detected.	
	6 Chloroethane	64.00	Compound Not Detected.	
	7 Trichlorofluoromethane	101.00	Compound Not Detected.	
	11 1,1-Dichloroethene	96.00	Compound Not Detected.	
	16 Methylene chloride	84.00	9.034 9.034 (0.657) 68000	5 5
,	19 trans-1,2-Dichloroethene	96.00	Compound Not Detected.	
	21 1,1-Dichloroethane	63.00	Compound Not Detected.	
	31 Chloroform	83.00	Compound Not Detected.	
	33 1,1,1-Trichloroethane	97.00	Compound Not Detected.	
	35 Carbon tetrachloride	117.00	Compound Not Detected.	
*	36 1,4-Difluorobenzene	114.00	16.037 16.038 (1.000) 2179707	50
	37 1,2-Dichloroethane	62.00	Compound Not Detected.	
	38 Benzene	78.00	Compound Not Detected.	
	39 Trichloroethene	95.00	Compound Not Detected.	
	41 1,2-Dichloropropane	63.00	Compound Not Detected.	
	43 Bromodichloromethane	83.00	Compound Not Detected.	
	45 2-Chloroethylvinyl ether	63.00	Compound Not Detected.	
	46 cis-1,3-Dichloropropene	75.00	Compound Not Detected.	
*	48 Chlorobenzene-d5	117.00	22.382 22.372 (1.000) 1718605	50
	49 Toluene	91.00	Compound Not Detected.	
	51 trans-1,3-Dichloropropene	75.00	Compound Not Detected.	
	52 1,1,2-Trichloroethane	83.00	Compound Not Detected.	
	55 Tetrachloroethene	166.00	Compound Not Detected.	
	56 Dibromochloromethane	129.00	Compound Not Detected.	
	58 Chlorobenzene	112.00	Compound Not Detected.	
	60 Ethylbenzene	91.00	Compound Not Detected.	

·\$ . . .

								CONCENTRA	TIONS	
			QUANT SIG					ON-COLUMN	FINAL	
Co	oqmo	unds	MASS	RT	EXP RT	REL RT	RESPONSE	( ug/L)	( ug/L)	
						*****				
	51	trans-1,3-Dichloropropene	75.00		19.874	Comp	ound Not De	etected.		
	52	1,1,2-Trichloroethane	83.00		20.169	Comp	oound Not De	etected.		
	55	Tetrachloroethene	166.00		20.779	Comp	ound Not De	etected.		
	56	Dibromochloromethane	129.00		21.162	Comp	ound Not De	etected.		
	58	Chlorobenzene	112.00		22.461	Comp	ound Not De	etected.		
	60	Ethylbenzene	91.00		22.608	Comp	oound Not De	etected.		
	66	Bromoform	173.00		24.330	Comp	ound Not De	etected.		
	68	1,1,2,2-Tetrachloroethane	83.00		24.753	Comp	ound Not De	etected.		
٠	69	1,4-Dichlorobenzene-d4	152.00	27.329	27.320	(1.000)	868353	50		
\$	91	Dibromofluoromethane	113.00	13.775	13.766	(1.001)	922258	51	51	
\$	92	1,2-Dichloroethane-d4	65.00	15.014	15.005	(0.936)	385525	46	46	
\$	93	Toluene-d8	98.00	19.283	19.274	(1.202)	1938486	48	48	
s	94	Bromofluorobenzene	174.00	24.900	24.890	(1.113)	810014	48	48	

Data File: /chem/ms5.i/a071896a.b/18Jul0801008.d Page 1

Report Date: 29-Jul-1996 16:04

### Quality Analytical Laboratories

Unknown Compounds Quantitation Report

Data file : /chem/ms5.i/a071896a.b/18Jul0801008.d

Client Smp ID: TRIP BLANK Lab Smp Id: MB370007

Inj Date: 18-JUL-1996 18:51
Operator: WLH/RLW
Smp Info: TRIP\_BLANK MB370007
Misc Info: 5MLS Autotune Date: 18-Jul-96 11:10:3

Inst ID: ms5.i

Comment

: /chem/ms5.i/a071896a.b/8260w5.m Method

Meth Date : 29-Jul-1996 15:28 ms5

Cal Date : 10-JUL-1996 22:29 Cal File: 10Jul2001020.d

Als bottle: 8

Target Version: 3.12

Dil Factor: 1.000 Integrator: HP RTE Compound Sublist: BROWN&ROOT.sub

Sample Matrix: WATER

Ouantitative Mode : Use RF of Nearest Std

- NO TENTATIVELY IDENTIFIED COMPOUNDS -

GC/MS SEMIVOLATILE ORGANICS

# CASE NARRATIVE GC/MS SEMIVOLATILE ORGANICS

QAL Lab	Reference No./SDG	MB370
Project:	: Brown & Root Coa	stal Systems Station

#### I. RECEIPT

No exceptions were encountered unless a Sample Receipt Exception Report is attached to the Chain-of-Custody included with this data package.

#### II. HOLDING TIMES

- A. Sample Preparation: All holding times were met.
- B. Sample Analysis: All holding times were met.

#### III. METHOD

Preparation: SW-846 3520A

Cleanup: N/A

Analysis: SW-846 8270A

#### IV. PREPARATION

Sample preparation proceeded normally.

#### V. ANALYSIS

- A. Calibration: All acceptance criteria were met.
- B. Blanks: All acceptance criteria were met.
- C. Surrogates: All acceptance criteria were met.
- D. Spikes: As requested, the matrix spikes were performed using a sample from sample delivery group MB370 (MB370003MS and MB370003MSD). Please note that the relative percent recoveries for 4-Nitrophenol were above the advisory QC criteria in the matrix spike and matrix spike duplicate. All other advisory criteria were met. A copy of the results is provided for your review.
- E. Samples: Sample analysis proceeded normally.
- F. Other: Please note that the Form 1's reflect the specified target list and that Azobenzene is reported as 1,2-Diphenylhydrazine.

A summary of the most current applicable method detection limits (MDLs) immediately follows the case narrative.

Lab Reference No./SDG: Page 2

I certify that this data package is in compliance with the terms and conditions agreed to by the client and QAL, Inc., both technically and for completeness except for the conditions noted above. Release of the data contained in this hardcopy data package has been authorized by the Laboratory Manager or designated person, as verified by the following signature.

Lab Reference No./SDG: \_\_\_\_
Page 3

CASE NARRATIVE
Addendum

# Sample Information

LAB SAMPLE ID	CLIENT SAMPLE ID	SAMPLE MATRIX	DATE SAMPLED	DATE EXTRACTED	DATE ANALYZED	SAMPLE pH <sup>1</sup>
						<u> </u>
MB370001	333-MW01-1	WATER	07/11/96	07/16/96	07/25/96	N/A
MB370002	333-MW02-1	WATER	07/11/96	07/16/96	07/26/96	N/A
MB370003	333-MW03-1	WATER	07/11/96	07/16/96	07/25/96	N/A
MB370003MS	333-MW03-1MS	WATER	07/11/96	07/16/96	07/25/96	N/A
MB370003MSD	333-MW03-1MD	WATER	07/11/96	07/16/96	07/25/96	N/A
MB370004	333-PC4-1	WATER	07/11/96	07/16/96	07/25/96	N/A
MB370005	333-MW01-1B	WATER	07/11/96	07/16/96	07/24/96	N/A
MB370006	333-MW03-1D	WATER	07/11/96	07/16/96	07/25/96	N/A
C07166B1	SBLK7S	WATER	N/A	07/16/96	07/24/96	N/A

<sup>1</sup> Applies to samples designated for purgeable VOA analysis only.

#### ORGANICS ANALYSIS METHOD DETECTION LIMITS

#### GC/MS SEMIVOLATILE ORGANICS

Laboratory	Name:	CH2M HILL	Sample Matrix:	WATER
Analytical	Method:	SW8270	Extraction Method:	SW3520
			D.	<b>D</b> L
			ug	J/L
	Acei	naphthene	0	.43
		naphthylene	0.	.41
•		tophenone		.47
		cetylaminofluorene		. 68
		minobiphenyl		.31
		line	-	.70
		hracene		.46
		mite		.28
		mite(DUP)		.54
		zidine		.78
		zo(a) anthracene		.34
		zo(a)pyrene		.31
		zo(b)fluoranthene		.68
•		zo(g,h,i)perylene		.72
		zo(k)fluoranthene		.65
		zoic acid		.32
		zyl alcohol		.51
		(2-Chloroethoxy) meth		.60
		(2-Chloroethyl) ether		.61
		(2-Ethylhexyl)phthal		.51
		romophenyl-phenyleth		.56
		ylbenzylphthalate bazole		.25
				.56 .28
		hloro-3-Methylphenol hloroaniline		.47
		hloronaphthalene		.68
		hloronaphthalene		.53
		hlorophenol		.78
		nicrophenoi hlorophenyl-phenylet		.35
		ysene ysene	_	.36
		n-butylphthalate		.24
		n-octylphthalate		.45
		enz(a,h)anthracene		.54
		enzofuran		.43
		-Dichlorobenzene		.75
		-Dichlorobenzene		.73
	•	-Dichlorobenzene		.79
		'-Dichlorobenzidine		.67
		-Dichlorophenol		.54
		-Dichlorophenol		.62
		thylphthalate		.61
		2-Dimethylbenz(a)an		.36
		'-Dimethylbenzidine		.96
	-,-		•	

#### ORGANICS ANALYSIS METHOD DETECTION LIMITS

#### GC/MS SEMIVOLATILE ORGANICS

Laboratory Name: CH2M HILL Sample Matrix: WATER Extraction Method: Analytical Method: SW8270 SW3520 MDL ug/L 2,4-Dimethylphenol 0.44 Dimethylphthalate 1.69 1.36 4,6-Dinitro-2-methylphenol 1,3-Dinitrobenzene 0.85 2,4-Dinitrophenol 1.91 2,4-Dinitrotoluene 0.54 0.53 2,6-Dinitrotoluene 1,2-Diphenylhydrazine 0.42 0.42 Ethyl methanesulfonate Fluoranthene 0.81 Fluorene 0.53 Hexachlorobenzene 1.10 Hexachlorobutadiene 0.90 Hexachlorocyclopentadiene 1.98 Hexachloroethane 0.79 131.38 Hexachlorophene Hexachloropropene 1.81 0.46 Indeno(1,2,3-cd)pyrene 0.61 Isophorone Isosafrole 0.39 6.24 Methapyrilene Methyl methanesulfonate 0.42 0.50 3-Methylcholanthrene 0.62 2-Methylnaphthalene 2-Methylphenol 0.42 3 & 4-Methylphenol 1.14 N-Nitroso-di-n-butylamine 0.41 N-Nitroso-di-n-propylamine 0.72 0.31 N-Nitrosodiethylamine N-Nitrosodimethylamine 0.60 N-Nitrosodiphenylamine (1) 0.95 N-Nitrosomethylethylamine 0.43 N-Nitrosomorpholine 0.65 N-Nitrosopiperidine 0.40 N-Nitrosopyrrolidine 0.48 0.68 Naphthalene 1,4-Naphthoquinone 123.50 1-Naphthylamine 1.38 2-Naphthylamine 0.81 5-Nitro-o-toluidine 0.64 2-Nitroaniline 1.34

3-Nitroaniline

4-Nitroaniline

1.03

1.39

#### ORGANICS ANALYSIS METHOD DETECTION LIMITS

#### GC/MS SEMIVOLATILE ORGANICS

Laboratory Name: CH2M HILL Sample Matrix: WATER

Analytical Method: SW8270 Extraction Method: SW3520

<del></del>	
	MDL
	ug/L
Nitrobenzene	0.57
2-Nitrophenol	1.23
4-Nitrophenol	1.44
4-Nitroquinoline-1-oxide	2.02
o-Toluidine	0.30
2,2'-Oxybis(1-chloropropane)	0.51
p-Dimethylaminoazobenzene	0.62
Pentachlorobenzene	0.54
Pentachloronitrobenzene	0.28
Pentachlorophenol	2.51
Phenacetin	0.83
Phenanthrene	0.42
Phenol	0.53
Phenyl-tert-butylamine	25.65
1,4-Phenylenediamine	23.31
2-Picoline	1.03
Pronamide	0.65
Pyrene	0.43
Pyridine	0.87
Safrole	0.50
1,2,4,5-Tetrachlorobenzene	0.79
2,3,4,6-Tetrachlorophenol	1.01
1,2,4-Trichlorobenzene	0.72
2,4,5-Trichlorophenol	0.70
2,4,6-Trichlorophenol	0.66
1,3,5-Trinitrobenzene	0.65

333-MW01-1

Q

Lab Name: CH2M HILL Contract: MB370

CAS NO.

Lab Code: MGM Case No.: MB370 SAS No.: SDG No.: MB370

Matrix: (soil/water) WATER Lab Sample ID: MB370001

Sample wt/vol: 1000 (g/mL) ML Lab File ID: H0006699.D

Level: (low/med) LOW Date Received: 07/13/96

% Moisture: not dec. dec. Date Extracted:07/16/96

Extraction: (SepF/Cont/Sonc) CONT Date Analyzed: 07/25/96

GPC Cleanup: (Y/N) N pH: 7.0 Dilution Factor: 1.0

COMPOUND

CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L

62-75-9-----N-Nitrosodimethylamine 10 108-95-2----Phenol 10 U 111-44-4-----bis(2-Chloroethyl)ether U 10 95-57-8-----2-Chlorophenol 10 U 541-73-1----1,3-Dichlorobenzene U 10 106-46-7-----1,4-Dichlorobenzene U 10 95-50-1----1,2-Dichlorobenzene 10 U 108-60-1----2,2'-oxybis(1-Chloroprop\_(1) 10 U 621-64-7----N-Nitroso-di-n-propylamine 10 U 67-72-1-----Hexachloroethane 10 U U 98-95-3-----Nitrobenzene 10 78-59-1-----Isophorone 10 U 88-75-5-----2-Nitrophenol 10 U 105-67-9-----2,4-Dimethylphenol 10 U 111-91-1-----bis (2-Chloroethoxy) methane 10 U 10 U 120-83-2----2,4-Dichlorophenol 120-82-1----1,2,4-Trichlorobenzene U 10 U 91-20-3-----Naphthalene 10 Ų 87-68-3-----Hexachlorobutadiene 10 59-50-7----4-Chloro-3-methylphenol 10 U 10 U 88-06-2----2,4,6-Trichlorophenol U 91-58-7----2-Chloronaphthalene 10 U 131-11-3-----Dimethylphthalate 10 U 606-20-2----2,6-Dinitrotoluene 10 208-96-8-----Acenaphthylene 10 U 83-32-9-----Acenaphthene 10 U 51-28-5----2,4-Dinitrophenol 50 U 50 U 100-02-7----4-Nitrophenol U 10 121-14-2----2,4-Dinitrotoluene 10 U 84-66-2-----Diethylphthalate\_\_ 10 U 86-73-7-----Fluorene 7005-72-3----4-Chlorophenyl-phenylether 10 U

(1) 2,2'-oxybis(1-Chloropropane) is known as bis(2-Chloroisopropyl) ether

FORM I SV-1

SW846 M

333-MW01-1

Q

10

10

10

10

U

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U

U

Lab Name: CH2M HILL Contract: MB370

COMPOUND

CAS NO.

Lab Code: MGM Case No.: MB370 SAS No.: SDG No.: MB370

Matrix: (soil/water) WATER Lab Sample ID: MB370001

Sample wt/vol: 1000 (g/mL) ML Lab File ID: H0006699.D

Level: (low/med) LOW Date Received: 07/13/96

% Moisture: not dec. dec. Date Extracted:07/16/96

Extraction: (SepF/Cont/Sonc) CONT Date Analyzed: 07/25/96

GPC Cleanup: (Y/N) N pH: 7.0 Dilution Factor: 1.0

CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L

534-52-1-----4,6-Dinitro-2-methylphenol 50 U 86-30-6----N-Nitrosodiphenylamine (3) 10 U 122-66-7----1,2-Diphenylhydrazine 10 U 101-55-3-----4-Bromophenyl-phenylether\_ U 10 118-74-1-----Hexachlorobenzene U 10 87-86-5-----Pentachlorophenol U 50 85-01-8-----Phenanthrene 10 U 120-12-7-----Anthracene 10 U 84-74-2----Di-n-butylphthalate 10 U U 206-44-0-----Fluoranthene 10 92-87-5-----Benzidine U 50 129-00-0-----Pyrene U 10 85-68-7-----Butylbenzylphthalate 10 U 56-55-3-----Benzo (a) anthracene 10 U 91-94-1-----3,3'-Dichlorobenzidine 20 U 218-01-9-----Chrysene 10 U 117-81-7-----bis(2-Ethylhexyl)phthalate U 10 117-84-0-----Di-n-octylphthalate U 10 205-99-2----Benzo (b) fluoranthene 10 U 207-08-9-----Benzo(k)fluoranthene 10 U

(3) - Cannot be separated from Diphenylamine

50-32-8-----Benzo (a) pyrene

193-39-5-----Indeno (1, 2, 3-cd) pyrene

53-70-3-----Dibenz(a,h)anthracene

191-24-2-----Benzo(g,h,i)perylene

FORM I SV-2



# SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: CH2M HILL Contract: MB370 333-MW01-1

Lab Code: MGM Case No.: MB370

SAS No.: SDG No.: MB370

Matrix: (soil/water) WATER

Lab Sample ID: MB370001

Sample wt/vol: 1000 (g/mL) ML

Lab File ID: H0006699.D

Level: (low/med) LOW

Date Received: 07/13/96

% Moisture: not dec. dec.

Date Extracted:07/16/96

Extraction: (SepF/Cont/Sonc) CONT

Date Analyzed: 07/25/96

GPC Cleanup: (Y/N) N pH: 7.0

Dilution Factor: 1.0

CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L

Number TICs found: 12

SW846

CONCENTRATIONS

Report Date: 05-Aug-1996 14:10

#### CH2M Hill Montgomery

SW846 SEMI-VOLATILES

Data file : /chem/ms3.i/h072596a.b/h0006699.d /

Lab Smp Id: MB370001 ✓ Client Smp ID: 333-MW01-1 ✓

Inj Date : 25-JUL-1996 13:51

Smp Info : MB370001 333-MW01-1

Misc Info :

Comment :

Method : /chem/ms3.i/h072596a.b/SV8270.m /

Meth Date: 05-Aug-1996 14:10 mjohnson Quant Type: ISTD

Cal Date : 25-JUL-1996 09:31 Cal File: h0006693.d

Als bottle: 8

Dil Factor: 1.000 / Integrator: HP RTE

Integrator: HP RTE Compound Sublist: brown&root.sub

Target Version: 3.10

_		-					CONCENTRA		TIONS	
			QUANT SIG				ON-C	COLUMN	FINAL	
Со	Compounds  2 N-Nitrosodimethylamine  \$ 6 2-Fluorophenol  \$ 9 Phenol-d5  10 Phenol  12 bis(2-Chloroethyl)ether  14 2-Chlorophenol  15 1,3-Dichlorobenzene  * 16 1,4-Dichlorobenzene-d4		MASS	RT	EXP RT REL RT	RESPONSE	(	NG)	( ug/L)	
==	===	22222222222222222	====	==	======	=======	===	=====	*****	
	2	N-Nitrosodimethylamine	74.00	Соп	mpound Not Detect	ed.				
\$	6	2-Fluorophenol	112.00	5.471	5.471 (0.811)	173974		100	52	
\$	9	Phenol-d5	99.00	6.284	6.291 (0.932)	234002		110	56	
	10	Phenol	94.00	Con	pound Not Detect	ed.				
	12	bis(2-Chloroethyl)ether	63.00	Соп	pound Not Detect	ed.				
	14	2-Chlorophenol	128.00	Con	pound Not Detect	ed.				
	15	1,3-Dichlorobenzene	146.00	Соп	pound Not Detect	ed.				
٠.	16	1,4-9ichlorobenzene-d4	152.00	6.746	6.753 (1.000)	41867		40	(Q)	
	17	1,4-Dichlorobenzene	146.00	Соп	mpound Not Detect	ed.				
	20	1,2-Dichlorobenzene	146.00	Соп	mpound Not Detect	ed.				
	22	2,2'-oxybis(1-Chloropropane)	45.00	Con	npound Not Detect	ed.				
	26	N-Nitroso-di-n-propylamine	70.00	Con	apound Not Detect	ed.			_	
	30	Hexachloroethane	117.00	Con	mpound Not Detect	ed.				
\$	31	Nitrobenzene-d5	82.00	7.457	7.464 (0.876)	89101		58	29	
	32	Nitrobenzene	77.00	Соп	mpound Not Detect	ed.				
	34	Isophorone	82.00	Con	mpound Not Detect	ed.			•	
	35	2-Nitrophenol	139.00	Con	mpound Not Detect	ed.				
	36	2,4-Dimethylphenol	107.00	Con	mpound Not Detect	ed.				
	38	bis(2-Chloroethoxy)methane	93.00	Con	mpound Not Detect	ed.				
~	39	2,4-Dichlorophenol	162.00	Con	mpound Not Detect	ed.				
	41	1,2,4-Trichlorobenzene	180.00	Con	mpound Not Detect	ed.				
*	42	Naphthalene-d8	136.00	8.512	8.519 (1.000)	153915		40		
	43	Naphthalene	128.00	Con	mpound Not Detect	ed.				
	47	<b>Hexachlorobutadiene</b>	225.00	Con	mpound Not Detect	ed.				
	50	4-Chloro-3-methylphenol	107.00	Соп	mpound Not Detect	ed.				
	55	2,4,6-Trichlorophenol	196.00	Con	mpound Not Detect	ed.				
\$	57	2-Fluorobiphenyl	172.00	10.439	10.453 (0.875)	130917		55	28	
	59	2-Chloronaphthalene	162.00	Con	mpound Not Detect	ed.				
	63	Dimethylphthalate	163.00	Cor	apound Not Detect	ed.				
								Å,	. 1	

Data File: /chem/ms3.i/h072596a.b/h0006699.d Report Date: 05-Aug-1996 14:10

			CONCENTRATIONS
	QUANT SIG		ON-COLUMN FINAL
Compounds	MASS	RT EXP RT REL RT RESPONS	E ( NG) ( ug/L)
=======================================	====	-=	= 2000000 0000000
65 2,6-Dinitrotoluene	165.00	Compound Not Detected.	
66 Acenaphthylene	152.00	Compound Not Detected.	
* 68 Acenaphthene-d10	164.00	11.934 11.941 (1.000) 8455	7 40
69 Acenaphthene	154.00	Compound Not Detected.	
70 2,4-Dinitrophenol	184.00	Compound Not Detected.	
71 4-Nitrophenol	65.00	Compound Not Detected.	
74 2,4-Dinitrotoluene	165.00	Compound Not Detected.	
78 Diethylphthalate	149.00	Compound Not Detected.	
79 fluorene	166.00	Compound Not Detected.	
80 4-Chlorophenyl-phenylether	204.00	Compound Not Detected.	
83 4,6-Dinitro-2-methylphenol	198.00	Compound Not Detected.	
84 N-Nitrosodiphenylamine (1)	169.00	Compound Not Detected.	
85 1,2-Diphenylhydrazine	77.00	Compound Not Detected.	
\$ 86 2,4,6-Tribromophenol	329.65	13.736 13.744 (1.151) 4436	57 110 54
89 4-Bromophenyl-phenylether	248.00	Compound Not Detected.	
90 Hexachlorobenzene	284.00	Compound Not Detected.	
92 Pentachlorophenol -	266.00	Compound Not Detected.	
* 95 Phenanthrene-d10	188.00	15.334 15.348 (1.000) 1218	9 40
96 Phenanthrene	178.00	Compound Not Detected.	
97 Anthracene	178.00	Compound Not Detected.	
99 Di-n-butylphthalate	149.00	Compound Not Detected.	
102 Fluoranthene	202.00	Compound Not Detected.	
103 Benzidine	184.00	Compound Not Detected.	
104 Pyrene	202.00	Compound Not Detected.	
\$ 106 Terphenyl-d14	244.00	19.195 19.203 (0.879) 13873	50 25
110 Butylbenzylphthalate	149.00	Compound Not Detected.	
, 112 Benzo(a)anthracene	228.00	Compound Not Detected.	
113 3,3'-Dichlorobenzidine	252.00	Compound Not Detected.	
* 114 Chrysene-d12	240.00	21.841 21.848 (1.000) 943	54 40
115 Chrysene	228.00	Compound Not Detected.	
116 bis(2-Ethylhexyl)phthalate	149.00	Compound Not Detected.	
117 Di-n-octylphthalate	149.00	Compound Not Detected.	
118 Benzo(b)fluoranthene	252.00	Compound Not Detected.	
120 Benzo(k)fluoranthene	252.00	Compound Not Detected.	
122 Benzo(a)pyrene	252.00	Compound Not Detected.	
* 123 Perylene-d12	264.00	25.314 25.329 (1.000) 918	01 40
125 Indeno(1,2,3-cd)pyrene	276.00	Compound Not Detected.	
126 Dibenz(a,h)anthracene	278.00	Compound Not Detected.	
127 Benzo(g,h,i)perylene	276.00	Compound Not Detected.	

QC Flag Legend

Q - Qualifier signal failed the ratio test.

Data File: /chem/ms3.i/h072596a.b/h0006699.d

Report Date: 05-Aug-1996 12:02

#### CH2M Hill Montgomery

Unknown Compounds Quantitation Report

Data file : /chem/ms3.i/h072596a.b/h0006699.d

Lab Smp Id: MB370001 Client Smp ID: 333-MW01-1

Inj Date : 25-JUL-1996 13:51

Smp Info : MB370001 333-MW01-1

Misc Info:

Comment :

Method : /chem/ms3.i/h072596a.b/SV8270.m

Meth Date: 05-Aug-1996 10:38 mjohnson

Als bottle: 8

Dil Factor: 1.000 Target Version: 3.10

Integrator: HP RTE Compound Sublist: brown&root.sub

Sample Matrix: WATER

Quantitative Mode : Use RF of Nearest Std

IS	STD	===	RT ====	AREA	AMOUNT
*	16	1,4-Dichlorobenzene-d4	6.746	257147	40.000
*	42	Naphthalene-d8	8.512	361913	40.000
*	68	Acenaphthene-d10	11.934	538876	40.000
*	95	Phenanthrene-d10	15.334	369531	40.000
*	114	Chrysene-d12	21.841	367428	40.000

		CONCENT	RATIONS		Q	JANT	
RT	AREA ON	-COL( NG)	FINAL( ug/L)	QUAL	LIBRARY	LIB ENTRY	CPND #
unknown	2-but	oxy eth	and 008	1619k cas	**************************************	232#24222	======
5.713	49750	8	4	0		0	16
Unknown				CAS	#:		
8.995	36324	4	2	0		0	42
Unknown				CAS	#:		
9.611	53225	6	3	0		0	42
Benzoic ad	cid, ethyl-			CAS	#: 28134-31	-8	
9.955	87180	10	5	55	NBS75K.l	9725	42

S . 5

Data File: /chem/ms3.i/h072596a.b/h0006699.d Report Date: 05-Aug-1996 12:02

	CONCENTRATIONS				QUANT			
RT	AREA	ON-COL( NG)	FINAL( ug/L)	QUAL	LIBRARY	LIB ENTRY	CPND #	
====	====		=========	====	======	=======	Z=====	
Unknown				CAS	· #:			
10.204	36576	4	2	0		0	42	
Unknown				CAS	S #:			
12.608	54130	4	2	0		0	68	
Unknown	•			CAS	s #:			
13.121	96702	7	4	0		0	68(L)	
Benzene, d	iethenyl-			CAS	s #: 1321-74-(	)		
14.176	38058	. 4	2	72	NBS75K.l	65326	95	
1-Naphthal	eneacetic	acid		CAS	s #: 86-87-3			
15.129	54142	6	3	95	NBS75K.l	19486	95	
Unknown				ÇAS	S #:			
16.345	63180	7	3	0		0	95	
Unknown Or	ganic Aci	d		CAS	s #:			
16.492	56541	6	3	0		0	95	
Unknown				CAS	S #:			
19.914	175750	19	10	0		0	114	

## QC Flag Legend

'L - Operator selected an alternate library search match.

Page 2

Data File: /chem/ms3.i/h072596a.b/h0006699.d Report Date: 31-Jul-1996 16:14

Page 3

Flag Legend

L - Operator selected an alternate library search match.

# SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

333-MW02-1

Lab Name: CH2M HILL

Contract: MB370

Lab Code: MGM Case No.: MB370

SAS No.:

SDG No.: MB370

Matrix: (soil/water) WATER

Lab Sample ID: MB370002

Sample wt/vol: 1000 (g/mL) ML

Lab File ID: H0006710.D

Level: (low/med) LOW

Date Received: 07/13/96

% Moisture: not dec. \_\_\_ dec.

Date Extracted:07/16/96

Extraction: (SepF/Cont/Sonc) CONT

Date Analyzed: 07/26/96

GPC Cleanup: (Y/N) N

pH: 7.0

Dilution Factor: 1.0

CONCENTRATION UNITS:

CAS NO.

COMPOUND

(ug/L or ug/Kg) UG/L

Q

62-75-9N-Nitrosodimethylamine		
108-95-2Phenol 111-44-4bis (2-Chloroethyl) ether 95-57-82-Chlorophenol 541-73-11, 3-Dichlorobenzene 106-46-71, 4-Dichlorobenzene 95-50-11, 2-Dichlorobenzene 108-60-12, 2'-oxybis (1-Chloroprop_(1)) 621-64-7	10	บ
95-57-82-Chlorophenol 541-73-11,3-Dichlorobenzene 106-46-71,4-Dichlorobenzene 95-50-11,2-Dichlorobenzene 108-60-12,2'-oxybis(1-Chloroprop_(1)) 621-64-7N-Nitroso-di-n-propylamine 67-72-1Hexachloroethane 98-95-3Nitrobenzene 78-59-1Isophorone 88-75-52-Nitrophenol 105-67-92,4-Dimethylphenol 111-91-1bis(2-Chloroethoxy)methane 120-83-22,4-Dichlorophenol 120-82-11,2,4-Trichlorobenzene	10	Ū
95-57-82-Chlorophenol 541-73-11,3-Dichlorobenzene 106-46-71,4-Dichlorobenzene 95-50-11,2-Dichlorobenzene 108-60-12,2'-oxybis(1-Chloroprop_(1)) 621-64-7N-Nitroso-di-n-propylamine 67-72-1Hexachloroethane 98-95-3Nitrobenzene 78-59-1Isophorone 88-75-52-Nitrophenol 105-67-92,4-Dimethylphenol 111-91-1bis(2-Chloroethoxy)methane 120-83-22,4-Dichlorophenol 120-82-11,2,4-Trichlorobenzene	10	Ū
541-73-11, 3-Dichlorobenzene         106-46-71, 4-Dichlorobenzene         95-50-11, 2-Dichlorobenzene         108-60-12, 2'-oxybis (1-Chloroprop_(1)         621-64-7Nitroso-di-n-propylamine         67-72-1Hexachloroethane         98-95-3Nitrobenzene         78-59-1Isophorone         88-75-52-Nitrophenol         105-67-92, 4-Dimethylphenol         111-91-1bis (2-Chloroethoxy) methane         120-83-22, 4-Dichlorophenol         120-82-11, 2, 4-Trichlorobenzene	10	Ü
106-46-71,4-Dichlorobenzene  95-50-11,2-Dichlorobenzene  108-60-12,2'-oxybis(1-Chloroprop_(1)) 621-64-7N-Nitroso-di-n-propylamine 67-72-1Hexachloroethane 98-95-3Nitrobenzene 78-59-1Isophorone 88-75-52-Nitrophenol 105-67-92,4-Dimethylphenol 111-91-1bis(2-Chloroethoxy)methane 120-83-22,4-Dichlorophenol 120-82-11,2,4-Trichlorobenzene	10	U
95-50-11,2-Dichlorobenzene  108-60-12,2'-oxybis(1-Chloroprop_(1) 621-64-7N-Nitroso-di-n-propylamine 67-72-1Hexachloroethane 98-95-3Nitrobenzene 78-59-1Isophorone 88-75-52-Nitrophenol 105-67-92,4-Dimethylphenol 111-91-1bis(2-Chloroethoxy)methane 120-83-22,4-Dichlorophenol 120-82-11,2,4-Trichlorobenzene	10	U
108-60-12,2'-oxybis(1-Chloroprop_(1)) 621-64-7N-Nitroso-di-n-propylamine_ 67-72-1Hexachloroethane_ 98-95-3Nitrobenzene_ 78-59-1Isophorone_ 88-75-52-Nitrophenol_ 105-67-92,4-Dimethylphenol_ 111-91-1bis(2-Chloroethoxy)methane_ 120-83-22,4-Dichlorophenol_ 120-82-11,2,4-Trichlorobenzene	10	U
621-64-7N-Nitroso-di-n-propylamine	10	Ū
67-72-1	10	Ŭ
98-95-3Nitrobenzene 78-59-1Isophorone 88-75-52-Nitrophenol 105-67-92,4-Dimethylphenol 111-91-1bis(2-Chloroethoxy)methane 120-83-22,4-Dichlorophenol 120-82-11,2,4-Trichlorobenzene	10	IJ
78-59-1Isophorone  88-75-52-Nitrophenol  105-67-92,4-Dimethylphenol  111-91-1bis(2-Chloroethoxy)methane  120-83-22,4-Dichlorophenol  120-82-11,2,4-Trichlorobenzene	10	Ū
88-75-52-Nitrophenol 105-67-92,4-Dimethylphenol 111-91-1bis(2-Chloroethoxy)methane 120-83-22,4-Dichlorophenol 120-82-11,2,4-Trichlorobenzene	10	Ū
105-67-92,4-Dimethylphenol 111-91-1bis(2-Chloroethoxy)methane 120-83-22,4-Dichlorophenol 120-82-11,2,4-Trichlorobenzene	10	Ū
111-91-1bis (2-Chloroethoxy) methane	10	IJ
120-83-22,4-Dichlorophenol	10	Ū
120-82-11,2,4-Trichlorobenzene	10	บ
	10	Ū
	10	Ū
87-68-3Hexachlorobutadiene	10	Ū
59-50-74-Chloro-3-methylphenol	10	Ü
88-06-22,4,6-Trichlorophenol	10	ט י
91-58-72-Chloronaphthalene	10	Ü
131-11-3Dimethylphthalate	10	Ü
606-20-22,6-Dinitrotoluene	10	Ü
208-96-8Acenaphthylene	10	Ü
83-32-9Acenaphthene	10	Ü
51-28-52,4-Dinitrophenol	50	บ
100-02-74-Nitrophenol	50	ΰ
121-14-22,4-Dinitrotoluene	10	ΰ
84-66-2Diethylphthalate	10	บ
86-73-7Fluorene	10	U
7005-72-34-Chlorophenyl-phenylether_	10	τ

(1) 2,2'-oxybis(1-Chloropropane) is known as bis(2-Chloroisopropyl) ether SW846 1778

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333-MW02-1

Lab Name: CH2M HILL Contract: MB370

Lab Code: MGM Case No.: MB370 SAS No.: SDG No.: MB370

Matrix: (soil/water) WATER Lab Sample ID: MB370002

Sample wt/vol: 1000 (g/mL) ML Lab File ID: H0006710.D

Level: (low/med) LOW Date Received: 07/13/96

% Moisture: not dec.\_\_\_\_\_ dec.\_\_\_ Date Extracted:07/16/96

Extraction: (SepF/Cont/Sonc) CONT Date Analyzed: 07/26/96

GPC Cleanup: (Y/N) N pH: 7.0 Dilution Factor: 1.0

CONCENTRATION UNITS: CAS NO. COMPOUND (ug/L or ug/Kg) UG/L Q 534-52-1----4,6-Dinitro-2-methylphenol 50 86-30-6----N-Nitrosodiphenylamine (3) 10 U 122-66-7----1,2-Diphenylhydrazine 10 U U 101-55-3----4-Bromophenyl-phenylether 10 U 118-74-1-----Hexachlorobenzene 10 U 87-86-5-----Pentachlorophenol 50 U 85-01-8-----Phenanthrene 10 120-12-7-----Anthracene 10 U 84-74-2----Di-n-butylphthalate 10 U 206-44-0-----Fluoranthene 10 U 92-87-5-----Benzidine 50 U U 129-00-0-----Pyrene 10 85-68-7-----Butylbenzylphthalate U 10 U 56-55-3-----Benzo (a) anthracene 10 91-94-1----3,3'-Dichlorobenzidine 20 U 218-01-9-----Chrysene 10 U 117-81-7-----bis(2-Ethylhexyl)phthalate U 10 U 117-84-0-----Di-n-octylphthalate 10 205-99-2----Benzo (b) fluoranthene U 10 Ū 207-08-9-----Benzo (k) fluoranthene 10 U 50-32-8-----Benzo (a) pyrene 10 U 193-39-5----Indeno(1,2,3-cd)pyrene 10 53-70-3-----Dibenz (a, h) anthracene 10 U U 191-24-2----Benzo(g,h,i)perylene 10

(3) - Cannot be separated from Diphenylamine FORM I SV-2

SW846 mil

# SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

333-MW02-1

Lab Name: CH2M HILL

Contract: MB370

Lab Code: MGM Case No.: MB370 SAS No.: SDG No.: MB370

Matrix: (soil/water) WATER

Lab Sample ID: MB370002

Sample wt/vol:

1000 (g/mL) ML

Lab File ID: H0006710.D

Level: (low/med)

LOW

Date Received: 07/13/96

% Moisture: not dec.\_\_\_\_ dec.

Date Extracted:07/16/96

Extraction: (SepF/Cont/Sonc) CONT

Date Analyzed: 07/26/96

GPC Cleanup: (Y/N) N pH: 7.0

Dilution Factor: 1.0

CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L

Number TICs found: 5

	,		·	
CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1. 123-42-2 2. 286-20-4 3. 111-76-2 4. 111-90-0 5. 6. 7. 8. 9. 10. 11. 12. 13. 14. 15. 16. 17. 18. 19. 20. 21. 22. 23.		5.176 5.396 5.711	EST. CONC.  2 5 3 4 12	· · ·
24. 25. 26. 27.				
27. 28. 29. 30.				

SW846

Report Date: 01-Aug-1996 13:06

# CH2M Hill Montgomery

## SW846 SEMI-VOLATILES

Data file : /chem/ms3.i/h072696a.b/h0006710.d —

Lab Smp Id: MB370002 / Client Smp ID: 333-MW02-1

Inj Date : 26-JUL-1996 10:56

Inst ID: ms3.i Operator : mjohnson

Smp Info : MB370002 333-MW301-1

Misc Info : Comment

Method : /chem/ms3.i/h072696a.b/SV8270.m /

Meth Date: 01-Aug-1996 13:06 mjohnson Quant Type: ISTD Cal Date: 26-JUL-1996 08:49 Cal File: h000670 Cal File: h0006708.d -

Als bottle: 4
Dil Factor: 1.000 —

Integrator: HP RTE Compound Sublist: brown&root.sub

Target Version: 3.10

						CONCENTR	RTIONS
		QUANT SIG				ON-COLUMN	FINAL
C	ompounds	MASS	RT	EXP RT REL RT	RESPONSE	( NG)	( ug/L)
=	******	====	==	=======================================	=======	======	======
	2 N-Nitrosodimethylamine	74.00	Соп	pound Not Detect	ed.		
\$	6 2-Fluorophenol	112.00	5.470	5.471 (0.811)	200928	140	72
	9 Phenoi-d5	99.00	6.283	6.292 (0.932)	247684	140	71
	10 Phenol	94.00	Соп	pound Not Detect	ed.		
	12 bis(2-Chloroethyl)ether	63.00	Соп	pound Not Detect	ed.		
	14 2-Chlorophenol	128.00	Com	pound Not Detect	ed.		
	15 1,3-Dichlorobenzene	146.00	Com	pound Not Detect	ed.		
*	16 1,4-Dichlorobenzene-d4	152.00	6.745	6.746 (1.000)	36526	40	(0)
	17 1,4-Dichlorobenzene	146.00	Com	pound Not Detect	ed.		
	20 1,2-Dichlorobenzene	146.00	Соп	pound Not Detect	ed.		
	22 2,2'-oxybis(1-Chloropropane)	45.00	Com	pound Not Detect	ed.		
	26 N-Nitroso-di-n-propylamine	70.00	Соп	pound Not Detect	ed.		
	30 Hexachloroethane	117.00	Соп	pound Not Detect	ed.		
\$	31 Nitrobenzene-d5	82.00	7.455	7.457 (0.876)	103324	81	41
	32 Nitrobenzene	77.00	Соп	pound Not Detect	ed.		
	34 Isophorone	82.00	Com	pound Not Detect	ed.		
	35 2-Nitrophenol	139.00	Соп	pound Not Detect	ed.		
	36 2,4-Dimethylphenol	107.00	Com	pound Not Detect	ed.		
•4	39 2,4-Dichlorophenol	162.00	Com	pound Not Detect	ed.		
	41 1,2,4-Trichlorobenzene	180.00	Com	pound Not Detect	ed.		
*	42 Naphthalene-d8	136.00	8.511	8.512 (1.000)	126800	40	
	43 Naphthalene	128.00	Соп	pound Not Detect	ed.		
	47 Hexachlorobutadiene	225.00	Com	pound Not Detect	ed.		
	50 4-Chloro-3-methylphenol	107.00	Com	pound Not Detect	ed.		
	55 2,4,6-Trichlorophenol	196.00	Com	pound Not Detect	ed.		
\$	57 2-Fluorobiphenyl	172.00	10.445	10.447 (0.875)	160811	84	42
	59 2-Chloronaphthalene	162.00	Com	pound Not Detect	ed.		
	63 Dimethylphthalate	163.00	Com	pound Not Detect	ed.		•
	65 2,6-Dinitrotoluene	165.00	Com	pound Not Detect	ed.	-2	÷ , t

		CONCENTRATIONS
	QUANT SIG	· ON-COLUMN FINAL
Compounds	MASS	RF EXP RT REL RT RESPONSE ( NG) ( ug/L)
=======================================	2222	00 20000 100000 2000000 1000000
66 Acenaphthylene	152.00	Compound Not Detected.
* 68 Acenaphthene-d10	164.00	11.933 11.934 (1.000) 67482 40
69 Acenaphthene	154.00	Compound Not Detected.
70 2,4-Dinitrophenol	184.00	Compound Not Detected.
71 4-Nitrophenol	65.00	Compound Not Detected.
74 2,4-Dinitrotoluene	165.00	Compound Not Detected.
78 Diethylphthalate	149.00	Compound Not Detected.
79 Fluorene	166.00	Compound Not Detected.
80 4-Chlorophenyl-phenyleth	ner 204.00	Compound Not Detected.
83 4,6-Dinitro-2-methylpher	nol 198.00	Compound Not Detected.
84 N-Nitrosodiphenylamine (	(1) 169.00	Compound Not Detected.
85 1,2-Diphenylhydrazine	77.00	Compound Not Detected.
\$ 86 2,4,6-Tribromophenol	329.65	13.735 13.737 (1.151) 49614 150 73
89 4-Bromophenyl-phenylethe	er 248.00	Compound Not Detected.
90 Hexachlorobenzene	284.00	Compound Not Detected.
92 Pentachlorophenol	266.00	Compound Not Detected.
* 95 Phenanthrene-d10	188.00	15.333 15.342 (1.000) 96707 40
96 Phenanthrene	178.00	Compound Not Detected.
97 Anthracene	178.00	Compound Not Detected.
99 Di-n-butylphthalate	149.00	Compound Not Detected.
102 Fluoranthene	202.00	Compound Not Detected.
103 Benzidine	184.00	Compound Not Detected.
104 Pyrene	202.00	Compound Not Detected.
\$ 106 Terphenyl-d14	244.00	19.194 19.196 (0.879) 166512 74 37
110 Butylbenzylphthalate	149.00	Compound Not Detected.
112 Benzo(a)anthracene	228.00	Compound Not Detected.
113 3,3'-Dichlorobenzidine	252.00	Compound Not Detected.
* * 114 Chrysene-d12	240.00	21.832 21.841 (1.000) 79701 40
115 Chrysene	228.00	Compound Not Detected.
116 bis(2-Ethylhexyl)phthala	ate 149.00	Compound Not Detected.
117 Di-n-octylphthalate	149.00	Compound Not Detected.
118 Benzo(b)fluoranthene	252.00	Compound Not Detected.
120 Benzo(k)fluoranthene	252.00	Compound Not Detected.
122 Benzo(a)pyrene	252.00	Compound Not Detected.
* 123 Perylene-d12	264.00	25.313 25.322 (1.000) 78846 40
125 Indeno(1,2,3-cd)pyrene	276.00	Compound Not Detected.
126 Dibenz(a,h)anthracene	278.00	Compound Not Detected.
127 Benzo(g,h,i)perylene	276.00	Compound Not Detected.

# QC Flag Legend

Q - Qualifier signal failed the ratio test.

Data File: /chem/ms3.i/h072696a.b/h0006710.d

Report Date: 01-Aug-1996 13:06

## CH2M Hill Montgomery

Unknown Compounds Quantitation Report

Data file : /chem/ms3.i/h072696a.b/h0006710.d

Lab Smp Id: MB370002 Client Smp ID: 333-MW02-1

Inj Date : 26-JUL-1996 10:56

Operator : mjohnson Smp Info : MB370002 333-MW301-1 Inst ID: ms3.i

Misc Info : Comment

Method : /chem/ms3.i/h072696a.b/SV8270.m

Meth Date: 01-Aug-1996 13:06 mjohnson

Cal Date : 26-JUL-1996 08:49 Cal File: h0006708.d

Als bottle: 4

Dil Factor: 1.000 Integrator: HP RTE Target Version: 3.10

Compound Sublist: brown&root.sub

Sample Matrix: WATER

Quantitative Mode : Use RF of Nearest Std

ISTD	RT	AREA	AMOUNT
========	====	=====	=====
* 16 1,4-Dichlorobenzene-d4	6.745	233683	40.000
* 114 Chrysene-d12	21.832	320869	40.000

			CONCENT	RATIONS		QU	JANT	
	RT	AREA	ON-COL( NG)	FINAL( ug/L)	QUAL	LIBRARY	LIB ENTRY	CPND #
•	====	. ====			====	======	=======	=====
	2-Pentanon	e, 4-hyd	roxy-4-methyl-		CAS	#: 123-42-2		
	5.176	23783	4	2	50	NBS75K.l	64275	16
	<del>Unknown -</del>				CAS	#: 286-21	0-4	
	5.396	61446	10	5	0		0	16
	Unknown At	<del>cohol.</del>			CAS	#: 111-76	-2-	
	5.711	31345	5	3	0		0	16
	Unkno⊌n At	<del>cohol •</del>	•		CAS	#: 111-90-	.0	
	6.510	44203	8	4	0		0	16
	Unknown				CAS	: <b>#:</b>		
	19.847	195935	24	12	0		0	114

\$ 15 m

333-MW03-1

H0006701.D

Lab File ID:

Lab Name: CH2M HILL Contract: MB370

Sample wt/vol:

Lab Code: MGM Case No.: MB370 SAS No.: SDG No.: MB370

1000 (g/mL) ML

Matrix: (soil/water) WATER

Lab Sample ID: MB370003

Level: (low/med) LOW Date Received: 07/13/96

% Moisture: not dec.\_\_\_\_ dec. Date Extracted:07/16/96

Extraction: (SepF/Cont/Sonc) CONT Date Analyzed: 07/25/96

GPC Cleanup: (Y/N) N pH: 7.0 Dilution Factor: 1.0

CONCENTRATION UNITS: CAS NO. COMPOUND (ug/L or ug/Kg) UG/L Q

62-75-9N-Nitrosodimethylamine	10	ט
108-95-2Phenol	10	Ū
111-44-4bis(2-Chloroethyl)ether	10	Ū
95-57-82-Chlorophenol	10	Ū
541-73-11,3-Dichlorobenzene	10	Ū
106-46-71,4-Dichlorobenzene	10	บั
95-50-11,2-Dichlorobenzene	10	. Ŭ
108-60-12,2'-oxybis(1-Chloroprop (1)	10	บั
621-64-7N-Nitroso-di-n-propylamine	10	Ü
67-72-1Hexachloroethane	10	บ
98-95-3Nitrobenzene	10	Ŭ
78-59-1Isophorone	10	บ
88-75-52-Nitrophenol	10	ט
105-67-92,4-Dimethylphenol	10	บ
111-91-1bis(2-Chloroethoxy) methane	10	บ
120-83-22,4-Dichlorophenol	10	บ
120-82-11,2,4-Trichlorobenzene	10	บ
91-20-3Naphthalene	10	บ
87-68-3Hexachlorobutadiene	10	Ü
59-50-74-Chloro-3-methylphenol	10	Ü
88-06-22,4,6-Trichlorophenol	10	, n
91-58-72-Chloronaphthalene	10	บ
131-11-3Dimethylphthalate	10	บ
606-20-22,6-Dinitrotoluene	10	Ü
208-96-8Acenaphthylene	10	Ü
83-32-9Acenaphthene	10	U U
51-28-52,4-Dinitrophenol	50	ט ו
100-02-74-Nitrophenol	50	ט
121-14-22,4-Dinitrotoluene	10	บ
84-66-2Diethylphthalate	10	U U
86-73-7Fluorene	10	บ
7005-72-34-Chlorophenyl-phenylether	10	וז ו
7005-72-34-Chitotophenyi-phenytecher	10	U
		l

(1) 2,2'-oxybis(1-Chloropropane) is known as bis(2-Chloroisopropyl) ether SW846 M FORM I SV-1

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333-MW03-1

Lab Name: CH2M HILL

Lab Code: MGM

Sample wt/vol:

Contract: MB370

Case No.: MB370 SAS No.: SDG No.: MB370

Matrix: (soil/water) WATER

Lab Sample ID: MB370003

1000 (g/mL) ML

Lab File ID: H0006701.D

Level: (low/med)

LOW

Date Received: 07/13/96

% Moisture: not dec. dec.

Date Extracted:07/16/96

Extraction: (SepF/Cont/Sonc) CONT

Date Analyzed: 07/25/96

GPC Cleanup: (Y/N) N

pH: 7.0

Dilution Factor: 1.0

CONCENTRATION UNITS:

CAS NO.

COMPOUND

(ug/L or ug/Kg) UG/L

Q

1		
534-52-14,6-Dinitro-2-methylphenol	50	ប
86-30-6N-Nitrosodiphenylamine (3)	10	บั
122-66-71,2-Diphenylhydrazine	10	Ū
101-55-34-Bromophenyl-phenylether	10	Ū
118-74-1Hexachlorobenzene	10	Ü
87-86-5Pentachlorophenol	50	Ū
85-01-8Phenanthrene	10	Ū
120-12-7Anthracene	10	Ū
84-74-2Di-n-butylphthalate	10	Ū
206-44-0Fluoranthene	10	ט
92-87-5Benzidine	50	Ū
129-00-0Pyrene	10	U
85-68-7Butylbenzylphthalate	10	U
56-55-3Benzo (a) anthracene	10	ប
91-94-13,3'-Dichlorobenzidine	20	Ū
218-01-9Chrysene	10	Ū
117-81-7bis(2-Ethylhexyl)phthalate	10	ប
117-84-0Di-n-octylphthalate	10	U
205-99-2Benzo (b) fluoranthene	10	U
207-08-9Benzo(k)fluoranthene	10	ט
50-32-8Benzo(a)pyrene	10	ั ซ
193-39-5Indeno (1, 2, 3-cd) pyrene	10	U
53-70-3Dibenz(a,h)anthracene	10	Ŭ
191-24-2Benzo(g,h,i)perylene	10	U

(3) - Cannot be separated from Diphenylamine

FORM I SV-2

ي پر څ

333-MW03-1

Lab Name: CH2M HILL

Contract: MB370

Lab Code: MGM

Case No.: MB370 SAS No.: SDG No.: MB370

Matrix: (soil/water) WATER

Lab Sample ID: MB370003

Sample wt/vol: 1000 (g/mL) ML

Lab File ID: H0006701.D

Level: (low/med)

Number TICs found: 6

Date Received: 07/13/96

% Moisture: not dec.\_\_\_\_ dec.\_\_\_

Date Extracted:07/16/96

Extraction: (SepF/Cont/Sonc) CONT

LOW

Date Analyzed: 07/25/96

GPC Cleanup: (Y/N) N

pH: 7.0

Dilution Factor: 1.0

CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L

	<u> </u>	1		
CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
=======================================	=======================================		==========	=====
1. 123-42-2	2-Pentanone, 4-hydroxy-4-met 7-Oxabicyclo[4.1.0] heptane	5.179	3	NJAB
2. 286-20-4	7-Oxabicyclo[4.1.0]heptane	5.398	3	VЛ
3. 111-76-2	Ethanol, 2-butoxy-	5.721	4	NJB
4.	Unknown	6.512	5	J
5.	Unknown	6.600	2	J
<u>6</u> .	Unknown	19.849	11	J
7				
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Report Date: 05-Aug-1996 14:11

# CH2M Hill Montgomery

SW846 SEMI-VOLATILES

Data file : /chem/ms3.i/h072596a.b/h0006701.d /

Lab Smp Id: MB370003 Client Smp ID: 333-MW03-1 /

Inj Date : 25-JUL-1996 15:10

Inst ID: ms3.i

Operator : mjohnson Smp Info : MB370003 333-MW03-1

Misc Info : Comment

Method : /chem/ms3.i/h072596a.b/SV8270.m / Meth Date : 05-Aug-1996 14:10 mjohnson Quant Type: ISTD

Cal File: h0006693.d / Cal Date : 25-JUL-1996 09:31

Als bottle: 10

Dil Factor: 1.000/ Integrator: HP RTE

Compound Sublist: brown&root.sub

Target Version: 3.10

	_		(	CONCENTRATIONS
		QUANT SIG	ON-	-COLUMN FINAL
Çc	mpounds	MASS	RT EXP RT REL RT RESPONSE (	NG) (ug/L)
==		====	== ====== ====== =====================	
	2 N-Nitrosodimethylamine	74.00	Compound Not Detected.	
\$	6 2-Fluorophenol	112.00	5.472 5.471 (0.811) 223654	160 81
\$	9 Phenol-d5	99.00	6.292 6.291 (0.933) 266947	160 78
	10 Phenol	94.00	Compound Not Detected.	
	12 bis(2-Chloroethyl)ether	63.00	Compound Not Detected.	
	14 2-Chlorophenol	128.00	Compound Not Detected.	
	15 1,3-Dichlorobenzene	146.00	Compound Not Detected.	
1 🛊	16 1,4-Dichlorobenzene-d4	152.00	6.747 6.753 (1.000) 34597	40 (Q)
	17 1,4-Dichlorobenzene	146.00	Compound Not Detected.	
	20 1,2-Dichlorobenzene	146.00	Compound Not Detected.	
	22 2,2'-oxybis(1-Chloropropane)	45.00	Compound Not Detected.	
	26 N-Nitroso-di-n-propylamine	70.00	Compound Not Detected.	
	30 Hexachloroethane	117.00	Compound Not Detected.	
\$	31 Nitrobenzene-d5	82.00	7.465 7.464 (0.877) 103013	81 40
	32 Nitrobenzene	77.00	Compound Not Detected.	
	34 Isophorone	82.00	Compound Not Detected.	
	35 2-Nitrophenol	139.00	Compound Not Detected.	
	36 2,4-Dimethylphenol	107.00	Compound Not Detected.	
	38 bis(2-Chloroethoxy)methane	93.00	Compound Not Detected.	
•	39 2,4-Dichlorophenol	162.00	Compound Not Detected.	
	41 1,2,4-Trichlorobenzene	180.00	Compound Not Detected.	
*	42 Naphthalene-d8	136.00	8.513 8.519 (1.000) 127803	40
	43 Naphthalene	128.00	Compound Not Detected.	
	47 Hexachlorobutadiene	225.00	Compound Not Detected.	
	50 4-Chloro-3-methylphenol	107.00	Compound Not Detected.	
	55 2,4,6-Trichlorophenol	196.00	Compound Not Detected.	
\$	57 2-Fluorobiphenyl	172.00	10.447 10.453 (0.875) 161158	78 39
	59 2-Chloronaphthalene	162.00	Compound Not Detected.	
	63 Dimethylphthalate	163.00	Compound Not Detected.	7
				÷ 3

pur mossion 000197

Data File: /chem/ms3.i/h072596a.b/h0006701.d Report Date: 05-Aug-1996 14:11

			CONCENTRATIONS
	QUANT SIG	. 0	N-COLUMN FINAL
Compounds	MASS	RT EXP RT REL RT RESPONSE	( NG) ( ug/L)
*********	====	22 20222 22222 222222	*******
65 2,6-Dinitrotoluene	165.00	Compound Not Detected.	
66 Acenaphthylene	152.00	Compound Not Detected.	
* 68 Acenaphthene-d10	164.00	11.942 11.941 (1.000) 73489	40
69 Acenaphthene	154.00	Compound Not Detected.	
70 2,4-Dinitrophenol	184.00	Compound Not Detected.	
71 4-Nitrophenol	65.00	Compound Not Detected.	
74 2,4-Dinitrotoluene	165.00	Compound Not Detected.	
.78 Diethylphthalate	149.00	Compound Not Detected.	
79 Fluorene	166.00	Compound Not Detected.	
80 4-Chlorophenyl-phenylether	204.00	Compound Not Detected.	
83 4,6-Dinitro-2-methylphenol	198.00	Compound Not Detected.	
84 N-Nitrosodiphenylamine (1)	169.00	Compound Not Detected.	
85 1,2-Diphenylhydrazine	77.00	Compound Not Detected.	
\$ 86 2,4,6-Tribromophenol	329.65	13.737 13.744 (1.150) 53473	150 75
89 4-Bromophenyl-phenylether	248.00	Compound Not Detected.	
90 Hexachlorobenzene	284.00	Compound Not Detected.	
92 Pentachlorophenol -	266.00	Compound Not Detected.	
* 95 Phenanthrene-d10 .	188,00	15.342 15.348 (1.000) 106536	40
96 Phenanthrene	178.00	Compound Not Detected.	
97 Anthracene	178.00	Compound Not Detected.	
99 Di-n-butylphthalate	149.00	Compound Not Detected.	
102 Fluoranthene	202.00	Compound Not Detected.	
103 Benzidine	184.00	Compound Not Detected.	
104 Pyrene	202.00	Compound Not Detected.	
\$ 106 Terphenyl-d14	244.00	19.204 19.203 (0.879) 191608	83 42
110 Butylbenzylphthalate	149.00	Compound Not Detected.	
112 Benzo(a)anthracene	228.00	Compound Not Detected.	
113 3,3'-Dichlorobenzidine	252.00	Compound Not Detected.	
* 114 Chrysene-d12	240.00	21.842 21.848 (1.000) 78248	40
115 Chrysene	228.00	Compound Not Detected.	
116 bis(2-Ethylhexyl)phthalate	149.00	Compound Not Detected.	
117 Di-n-octylphthalate	149.00	Compound Not Detected.	
118 Benzo(b)fluoranthene	252.00	Compound Not Detected.	
120 Benzo(k)fluoranthene	252.00	Compound Not Detected.	
122 Benzo(a)pyrene	252.00	Compound Not Detected.	
* 123 Perylene-d12	264.00	25.322 25.329 (1.000) 67484	40
125 Indeno(1,2,3-cd)pyrene	276.00	Compound Not Detected.	
126 Dibenz(a,h)anthracene	278.00	Compound Not Detected.	
127 Benzo(g,h,i)perylene	276.00	Compound Not Detected.	

# QC Flag Legend

Q - Qualifier signal failed the ratio test.

Page 3

Report Date: 31-Jul-1996 16:15

# CH2M Hill Montgomery

Unknown Compounds Quantitation Report

Data file : /chem/ms3.i/h072596a.b/h0006701.d

Lab Smp Id: MB370003 Client Smp ID: 333-MW03-1

Inj Date : 25-JUL-1996 15:10

Smp Info : MB370003 333-MW03-1

Misc Info :

Comment :

Method : /chem/ms3.i/h072596a.b/SV8270.m

Meth Date: 31-Jul-1996 16:11 mjohnson

Als bottle: 10

Dil Factor: 1.000 Target Version: 3.10

Integrator: HP RTE Compound Sublist: brown&root.sub

Sample Matrix: WATER

Quantitative Mode : Use RF of Nearest Std

ISTD	RT	AREA	TRUOMA
*****	====	======	=====
* 16 1,4-Dichlorobenzene-d4	6.747	222144	40.000
* 114 Chrysene-d12	21.842	323072	40.000

		CONCE	NTRATIONS		QU	ANT	
RT	AREA	ON-COL( NG	) FINAL( ug/L)	QUAL	LIBRARY	LIB ENTRY	CPND #
, ==== -	. ====	==========	= ==========	====	======	22222222	=====
2-Pentanon	e, 4-hyd	roxy-4-methyl		- CAS	s #: 123-42-2		
5.179	<b>3</b> 5560	6	3	56	NBS75K.l	64275	16
→ Unknown	•			CAS	; #: 286-z	0-4	
5.398	33343	6	3	0		0	16
Unknown Al	cohol			CAS	s #: 111-76	- 2	
5.721	45884	8	4	0		0	16
Unknown At	conol			CAS	S #:		
6.512	53627	10	5	0		0	16
Unknown At	cohol			CAS	S #:		
6.600	22722	4	2	0		0	16
Unknown				CAS	s <b>#</b> :		
19.849	175054	22	11	0		0	114

EPA SAMPLE NO.

333-PC4-1

Lab Name: CH2M HILL

Contract: MB370

Lab Code: MGM

Case No.: MB370

SAS No.:

SDG No.: MB370

Matrix: (soil/water) WATER

Lab Sample ID: MB370004

Sample wt/vol:

Lab File ID:

H0006702.D

Level: (low/med) LOW

1000 (g/mL) ML

Date Received: 07/13/96

% Moisture: not dec. dec.

Date Extracted: 07/16/96

Extraction: (SepF/Cont/Sonc) CONT

Date Analyzed: 07/25/96

GPC Cleanup: (Y/N) N

pH: 7.0

Dilution Factor: 1.0

CONCENTRATION UNITS:

CAS NO.

COMPOUND

(ug/L or ug/Kg) UG/L

Q

108-95-2Phenol 111-44-4bis (2-Chloroethyl) ether 95-57-82-Chlorophenol 10	<del></del>	· <del></del> -	
108-95-2Phenol       10       II         111-44-4	62-75-9N-Nitrosodimethylamine	10	ប
111-44-4		10	บ
95-57-82-Chlorophenol 10	111-44-4bis(2-Chloroethyl)ether		Ü
541-73-11,3-Dichlorobenzene       10         106-46-71,4-Dichlorobenzene       10         95-50-11,2-Dichlorobenzene       10         108-60-12,2'-oxybis (1-Chloroprop_(1))       10         621-64-7Nitroso-di-n-propylamine       10         67-72-1Hexachloroethane       10         98-95-3Nitrobenzene       10         78-59-1Isophorone       10         88-75-52-Nitrophenol       10         105-67-92,4-Dimethylphenol       10         111-91-1bis (2-Chloroethoxy) methane       10         120-83-22,4-Dichlorophenol       10         120-82-11,2,4-Trichlorobenzene       10         91-20-3Naphthalene       10         87-68-3Naphthalene       10         87-68-3		i	σ
106-46-71,4-Dichlorobenzene       10         95-50-11,2-Dichlorobenzene       10         108-60-12,2'-oxybis(1-Chloroprop_(1))       10         621-64-7N-Nitroso-di-n-propylamine       10         67-72-1Hexachloroethane       10         98-95-3Nitrobenzene       10         78-59-1Isophorone       10         88-75-52-Nitrophenol       10         105-67-92,4-Dimethylphenol       10         111-91-1bis(2-Chloroethoxy)methane       10         120-83-22,4-Dichlorophenol       10         120-83-22,4-Dichlorophenol       10         120-83-1		10	U
95-50-1			Ü
108-60-12,2'-oxybis(1-Chloroprop_(1))       10         621-64-7N-Nitroso-di-n-propylamine       10         67-72-1Hexachloroethane       10         98-95-3Nitrobenzene       10         78-59-1Isophorone       10         88-75-52-Nitrophenol       10         105-67-92,4-Dimethylphenol       10         111-91-1		1	Ū
621-64-7N-Nitroso-di-n-propylamine 10		10	บ
67-72-1	621-64-7Nitroso-di-n-propylamine		Ŭ
98-95-3Nitrobenzene       10       10         78-59-1Isophorone       10       10         88-75-52-Nitrophenol       10       10         105-67-92,4-Dimethylphenol       10       10         111-91-1bis (2-Chloroethoxy) methane       10       10         120-83-22,4-Dichlorophenol       10       10         120-82-11,2,4-Trichlorobenzene       10       10         91-20-3Naphthalene       10       10         87-68-3	67-72-1Hexachloroethane		บั
78-59-1			Ŭ
88-75-52-Nitrophenol       10         105-67-92,4-Dimethylphenol       10         111-91-1bis(2-Chloroethoxy)methane       10         120-83-22,4-Dichlorophenol       10         120-82-11,2,4-Trichlorobenzene       10         91-20-3Naphthalene       10         87-68-3Hexachlorobutadiene       10         59-50-74-Chloro-3-methylphenol       10         88-06-22,4,6-Trichlorophenol       10         91-58-72-Chloronaphthalene       10         131-11-3Dimethylphthalate       10         606-20-22,6-Dinitrotoluene       10         208-96-8Acenaphthene       10         51-28-52,4-Dinitrophenol       50         100-02-74-Nitrophenol       50         121-14-22,4-Dinitrotoluene       10         84-66-2Diethylphthalate       10         86-73-7Fluorene       10			บ
105-67-92,4-Dimethylphenol       10         111-91-1bis(2-Chloroethoxy)methane       10         120-83-22,4-Dichlorophenol       10         120-82-11,2,4-Trichlorobenzene       10         91-20-3Naphthalene       10         87-68-3Naphthalene       10         59-50-7			บ
111-91-1bis (2-Chloroethoxy) methane       10       10         120-83-22,4-Dichlorophenol       10       10         120-82-11,2,4-Trichlorobenzene       10       10         91-20-3Naphthalene       10       10         87-68-3Hexachlorobutadiene       10       10         59-50-74-Chloro-3-methylphenol       10       10         88-06-22,4,6-Trichlorophenol       10       10         91-58-72-Chloronaphthalene       10       10         131-11-3Dimethylphthalate       10       10         606-20-22,6-Dinitrotoluene       10       10         208-96-8Acenaphthylene       10       10         83-32-9Acenaphthene       10       10         51-28-52,4-Dinitrophenol       50       10         100-02-74-Nitrophenol       50       10         121-14-22,4-Dinitrotoluene       10       10         84-66-2Fluorene       10       10			บั
120-83-22,4-Dichlorophenol       10         120-82-11,2,4-Trichlorobenzene       10         91-20-3Naphthalene       10         87-68-3Hexachlorobutadiene       10         59-50-74-Chloro-3-methylphenol       10         88-06-22,4,6-Trichlorophenol       10         91-58-72-Chloronaphthalene       10         131-11-3Dimethylphthalate       10         606-20-22,6-Dinitrotoluene       10         208-96-8Acenaphthylene       10         83-32-9Acenaphthene       10         100-02-74-Nitrophenol       50         121-14-22,4-Dinitrotoluene       10         84-66-2Diethylphthalate       10         86-73-7Fluorene       10			Ŭ
120-82-11,2,4-Trichlorobenzene       10       0         91-20-3Naphthalene       10       0         87-68-3			บั
91-20-3Naphthalene       10       10         87-68-3			์ บั
87-68-3			Ū
59-50-74-Chloro-3-methylphenol       10       10         88-06-22,4,6-Trichlorophenol       10       10         91-58-72-Chloronaphthalene       10       10         131-11-3Dimethylphthalate       10       10         606-20-22,6-Dinitrotoluene       10       10         208-96-8Acenaphthylene       10       10         83-32-9Acenaphthene       10       10         51-28-52,4-Dinitrophenol       50       10         100-02-74-Nitrophenol       50       10         121-14-22,4-Dinitrotoluene       10       10         84-66-2Diethylphthalate       10       10         86-73-7			ט
88-06-22,4,6-Trichlorophenol       10         91-58-72-Chloronaphthalene       10         131-11-3Dimethylphthalate       10         606-20-22,6-Dinitrotoluene       10         208-96-8Acenaphthylene       10         83-32-9Acenaphthene       10         51-28-52,4-Dinitrophenol       50         100-02-74-Nitrophenol       50         121-14-22,4-Dinitrotoluene       10         84-66-2Diethylphthalate       10         86-73-7			บ็
91-58-72-Chloronaphthalene       10         131-11-3Dimethylphthalate       10         606-20-22,6-Dinitrotoluene       10         208-96-8Acenaphthylene       10         83-32-9Acenaphthene       10         51-28-52,4-Dinitrophenol       50         100-02-74-Nitrophenol       50         121-14-22,4-Dinitrotoluene       10         84-66-2Diethylphthalate       10         86-73-7Fluorene       10			
131-11-3Dimethylphthalate       10       0         606-20-22,6-Dinitrotoluene       10       0         208-96-8Acenaphthylene       10       0         83-32-9Acenaphthene       10       0         51-28-52,4-Dinitrophenol       50       0         100-02-74-Nitrophenol       50       0         121-14-22,4-Dinitrotoluene       10       0         84-66-2Diethylphthalate       10       0         86-73-7			บ
606-20-22,6-Dinitrotoluene       10         208-96-8Acenaphthylene       10         83-32-9Acenaphthene       10         51-28-52,4-Dinitrophenol       50         100-02-74-Nitrophenol       50         121-14-22,4-Dinitrotoluene       10         84-66-2Diethylphthalate       10         86-73-7Fluorene       10			Ιŭ
208-96-8			Ŭ
83-32-9			บ็
51-28-52,4-Dinitrophenol       50         100-02-74-Nitrophenol       50         121-14-22,4-Dinitrotoluene       10         84-66-2Diethylphthalate       10         86-73-7Fluorene       10			υ
100-02-74-Nitrophenol       50       U         121-14-22,4-Dinitrotoluene       10       U         84-66-2Diethylphthalate       10       U         86-73-7Fluorene       10       U			ไ บั
121-14-22,4-Dinitrotoluene       10       0         84-66-2Diethylphthalate       10       0         86-73-7Fluorene       10       0			ដ
84-66-2Diethylphthalate 10 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0			ľ
86-73-7Fluorene10			ט
			ľ
1002-15-2			"
į l	7003-72 34-Cittotophenyt-phenytechet		1

(1) 2,2'-oxybis(1-Chloropropane) is known as bis(2-Chloroisopropyl) ether SW846 1971

FORM I SV-1

### 1C SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

333-PC4-1

Lab Name: CH2M HILL Contract: MB370

Lab Code: MGM Case No.: MB370 SAS No.: SDG No.: MB370

Matrix: (soil/water) WATER Lab Sample ID: MB370004

Sample wt/vol: 1000 (g/mL) ML Lab File ID: H0006702.D

Level: (low/med) LOW Date Received: 07/13/96

% Moisture: not dec. Date Extracted:07/16/96

Extraction: (SepF/Cont/Sonc) CONT Date Analyzed: 07/25/96

GPC Cleanup: (Y/N) N pH: 7.0 Dilution Factor: 1.0

CONCENTRATION UNITS:

CAS NO. COMPOUND (ug/L or ug/Kg) UG/L Q

(3) - Cannot be separated from Diphenylamine FORM I SV-2

SW846 YMM

333-PC4-1

Lab Name: CH2M HILL

Contract: MB370

Lab Code: MGM

Case No.: MB370 SAS No.: SDG No.: MB370

Matrix: (soil/water) WATER

Lab Sample ID: MB370004

Sample wt/vol: 1000 (g/mL) ML

Lab File ID: H0006702.D

Level: (low/med) LOW

Date Received: 07/13/96

% Moisture: not dec.\_\_\_\_ dec.

Date Extracted:07/16/96

Extraction: (SepF/Cont/Sonc) CONT

Date Analyzed: 07/25/96

GPC Cleanup: (Y/N) N

pH: 7.0

Dilution Factor: 1.0

CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L

Number TICs found: 13

1	T		<del></del>	<del></del> 1
CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1 102 42 2	2 Dantanana A hadaaaa A mah		========	=====
1. 123-42-2	2-Pentanone, 4-hydroxy-4-met		5	NJAB
2. 286-20-4	7-Oxabicyclo[4.1.0]heptane	5.406	7	NJ
3. 822-67-3	2-Cyclohexen-1-ol	5.626	2	ŊJ
4. 111-76-2	Ethanol, 2-butoxy-	5.721	4	NJB
5. 930-68-7	2-Cyclohexen-1-one	6.044	2	. NJ
6. 111-90-0	Ethanol, 2-(2-ethoxyethoxy)-		4	ПЛ
7. 13429-07-7	2-Propanol, 1-(2-methoxyprop	6.600	2	NJ
8. 2039-89-6	Benzene, 2-ethenyl-1,4-dimet		2	NJ
9. 544-63-8	Tetradecanoic acid	14.426	4	ŊJ
10.	Unknown Hydrocarbon	16.339	16	J
11.	Unknown Hydrocarbon	16.405	7	J
12. 57-10-3	Hexadecanoic acid	16.493	4	IJ
13.	Unknown	19.864	14	J
14				' I
1 15.				
16.				
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Sec. 3.

Report Date: 05-Aug-1996 14:11

# CH2M Hill Montgomery

## SW846 SEMI-VOLATILES

Data file : /chem/ms3/i/h072596a.b/h0006702.d/

Client Smp ID: 333-PC4-1 Lab Smp Id: MB370004

Inj Date : 25-JUL-1996 15:50

Operator : mjohnson Inst ID: ms3.i

Smp Info : MB370004 333-PC4-1

Misc Info :

Comment Method : /chem/ms3.i/h072596a.b/SV8270.m /

Meth Date: 05-Aug-1996 14:10 mjohnson Quant Type: ISTD

Cal Date : 25-JUL-1996 09:31 Cal File: h0006693.d /

Als bottle: 11

Dil Factor: 1.000 / Integrator: HP RTE Compound Sublist: brown&root.sub

Target Version: 3.10

-					CONCENTR	ATIONS
	QUANT SIG				ON-COLUMN	FINAL
Compounds	MASS	RT	EXP RT REL RT	RESPONSE	( NG)	( ug/L)
#######################################	====	==	=======================================	*======	******	*****
2 N-Nitrosodimethylamine	74.00	Com	pound Not Detect	ed.		
\$ 6 2-Fluorophenol	112.00	5.472	5.471 (0.811)	240168	180	91
\$ 9 Phenol-d5	99.00	6.293	6.291 (0.933)	298463	180	90
10 Phenol	94.00	Com	pound Not Detect	ed.		
12 bis(2-Chloroethyl)ether	63.00	Com	pound Not Detect	ed.		
14 2-Chlorophenol	128.00	Com	pound Not Detect	ed.		
15 1,3-Dichlorobenzene	146.00	Com	pound Not Detect	ed.		
* 16 1,4-Dichlorobenzene-d4	152.00	6.747	6.753 (1.000)	33303	40	(Q)
17 1,4-Dichlorobenzene	146.00	Com	pound Not Detect	ed.		
20 1,2-Dichlorobenzene	146.00	· Com	pound Not Detect	ed.		
22 2,2'-oxybis(1-Chloropropane)	45.00	Com	pound Not Detect	ed.		
26 N-Nitroso-di-n-propylamine	70.00	Com	pound Not Detect	ed.		
30 Hexachloroethane	117.00	Con	pound Not Detect	ed.		
\$ 31 Nitrobenzene-d5	82.00	7.465	7.464 (0.877)	119064	93	46
32 Nitrobenzene	77.00	Соп	pound Not Detect	ed.		
34 Isophorone	82.00	Соп	pound Not Detect	ed.		
35 2-Nitrophenol	139.00	Con	pound Not Detect	ed.		
36 2,4-Dimethylphenol	107.00	Соп	pound Not Detect	ed.		
38 bis(2-Chloroethoxy)methane	93.00	Соп	pound Not Detect	ed.		
39 2,4-Dichlorophenol	162.00	Соп	pound Not Detect	ed.		
41 1,2,4-Trichlorobenzene	180.00	Соп	pound Not Detect	ed.		
* 42 Naphthalene-d8	136.00	8,513	8.519 (1.000)	129434	40	
43 Naphthalene	128.00	Con	pound Not Detect	ed.		
47 Hexachlorobutadiene	225.00	Соп	pound Not Detect	ed.		
50 4-Chloro-3-methylphenol	107.00	Соп	pound Not Detect	ed.		
55 2,4,6-Trichlorophenol	196.00	Соп	pound Not Detect	ed.		
\$ 57 2-Fluorobiphenyl	172.00	10.448	10.453 (0.875)	181745	86	43
59 2-Chloronaphthalene	162.00	Соп	pound Not Detect	ed.		
63 Dimethylphthalate	163.00	. Соп	pound Not Detect	ed.	•	
					. 2	, }

QUANT S  Compounds	G ON-COLUMN FINAL  RT EXP RT REL RT RESPONSE ( NG) ( ug/L)  == =================================
65 2,6-Dinitrotoluene 165.00 66 Acenaphthylene 152.00	Compound Not Detected. Compound Not Detected. 11.942 11.941 (1.000) 75498 40 Compound Not Detected.
65 2,6-Dinitrotoluene 165.00 66 Acenaphthylene 152.00	Compound Not Detected. Compound Not Detected. 11.942 11.941 (1.000) 75498 40 Compound Not Detected.
66 Acenaphthylene 152.00	Compound Not Detected. 11.942 11.941 (1.000) 75498 40 Compound Not Detected.
	11.942 11.941 (1.000) 75498 40 Compound Not Detected.
* 68 Acenaphthene-d10 164.00	Compound Not Detected.
	•
69 Acenaphthene 154.00	Compound Not Detected.
70 2,4-Dinitrophenol 184.00	
71 4-Nitrophenol 65.00	Compound Not Detected.
74 2,4-Dinitrotoluene 165.00	Compound Not Detected.
78 Diethylphthalate 149.00	Compound Not Detected.
79 Fluorene 166.00	Compound Not Detected.
80 4-Chlorophenyl-phenylether 204.00	Compound Not Detected.
83 4,6-Dinitro-2-methylphenol 198.00	Compound Not Detected.
84 N-Nitrosodiphenylamine (1) 169.00	Compound Not Detected.
85 1,2-Diphenylhydrazine 77.00	Compound Not Detected.
\$ 86 2,4,6-Tribromophenol 329.65	13.745 13.744 (1.151) 58257 160 79
89 4-Bromophenyl-phenylether 248.00	Compound Not Detected.
90 Hexachlorobenzene 284.00	Compound Not Detected.
92 Pentachlorophenol - 266.00	Compound Not Detected.
* 95 Phenanthrene-d10 · 188.00	15.342 15.348 (1.000) 107731 40
96 Phenanthrene 178.00	Compound Not Detected.
97 Anthracene 178.00	Compound Not Detected.
99 Di-n-butylphthalate 149.00	Compound Not Detected.
102 Fluoranthene 202.00	Compound Not Detected.
103 Benzidine 184.00	Compound Not Detected.
704 Pyrene 202.00	Compound Not Detected.
\$ 106 Terphenyl-d14 244.00	19.204 19.203 (0.879) 188790 80 40
110 Butylbenzylphthalate 149.00	Compound Not Detected.
112 Benzo(a)anthracene 228.00	Compound Not Detected.
113 3,3' Dichlorobenzidine 252.00	Compound Not Detected.
* 114 Chrysene-d12 240.00	21.842 21.848 (1.000) 79930 40
115 Chrysene 228.00	Compound Not Detected.
116 bis(2-Ethylhexyl)phthalate 149.00	Compound Not Detected.
117 Di-n-octylphthalate 149.00	Compound Not Detected.
118 Benzo(b)fluoranthene 252.00	Compound Not Detected.
120 Benzo(k)fluoranthene 252.00	Compound Not Detected.
122 Benzo(a)pyrene 252.00	Compound Not Detected.
* 123 Perylene-d12 264.00	25.323 25.329 (1.000) 71160 40
125 Indeno(1,2,3-cd)pyrene 276.00	Compound Not Detected.
126 Dibenz(a,h)anthracene 278.00	Compound Not Detected.
127 Benzo(g,h,i)perylene 276.00	Compound Not Detected.

# QC Flag Legend

Q - Qualifier signal failed the ratio test.

Data File: /chem/ms3.i/h072596a.b/h0006702.d

Report Date: 31-Jul-1996 16:15

Page 3

# CH2M Hill Montgomery

Unknown Compounds Quantitation Report

Data file : /chem/ms3.i/h072596a.b/h0006702.d

Lab Smp Id: MB370004 Client Smp ID: 333-PC4-1

Inj Date : 25-JUL-1996 15:50

Operator : mjohnson Inst ID: ms3.i

Smp Info : MB370004 333-PC4-1

Misc Info : Comment

Method : /chem/ms3.i/h072596a.b/SV8270.m

Meth Date : 31-Jul-1996 16:11 mjohnson Cal Date : 25-JUL-1996 09:31 Cal File: h0006693.d

Als bottle: 11

Dil Factor: 1.000 Target Version: 3.10

Integrator: HP RTE Compound Sublist: brown&root.sub

Sample Matrix: WATER

Quantitative Mode : Use RF of Nearest Std

IS	STD	RT ====	AREA	AMOUNT
*	16 1,4-Dichlorobenzene-d4	6.747	214519	40.000
*	42 Naphthalene-d8	8.513	320357	40.000
*	95 Phenanthrene-d10	15.342	343588	40.000
*	114 Chrysene-d12	21.842	308127	40.000

		C	ONCENT	RATIONS		QI	JANT	
RT	AREA	ON-COL(	NG)	FINAL( ug/L)	QUAL	LIBRARY	LIB ENTRY	CPND #
====	.====	=======	====	*******	====	*=====	*******	=====
/2-Pentanone,	4-hyd	roxy-4-me	thvl-		CAS	; #: 123-42 <b>-</b> 2		
5.179	55946		10	5	64	NBS75K.l	64275	16
-Unknown					CAS	#: 286 - 2	0-4	
5.406	74977	•	14	7	0		0	16
~ 2-Cyclohexen	1-ol				CAS	; #: 822-67-3		
5.626	24130		4	2	64	NBS75K.l	63202	16
Unknown Alec	hol				CAS	: #: III - 76	٠ عـ	
5.721	48563		9	4	0		0	16

Data File: /chem/ms3.i/h072596a.b/h0006702.d Report Date: 31-Jul-1996 16:15

Page 4

		CON	CENTRA	TIONS		QU	IANT	
RT	AREA	ON-COL(	NG)	FINAL( ug/L)	QUAL	LIBRARY	LIB ENTRY	CPND #
====	====	========			====	======	========	*****
2-Cycloh	exen-1-one				CA	s #: 930-68-7		
6.044	22685	4		2	90	NBS75K.l	1088	16
-Unknown 7	<del>tcoho</del> l				CA	s #: 111-90-	6	
6.513	40035	7	•	4	0		0	16
_Linknoun-	Ucohol				CA	S#: 13429.	- 07-7	
6.600	22753	4		2	0		0	16
Benzene,	2-ethenyl	-1,4-dimeth	yl-		CA	s #: 2039-89-6	5	
8.110		. 4		2	96	NBS75K.l	5879	42
Tetradec	anoic acid				CA	s #: 544-63-8		
	61359	•	,	4	97	NBS75K.l	70843	95
Unknown	hydroc	arbon			CA	s #:		
16.339	•	33		16	0		0	95
Unknown	hydroc	carbon	-		CA	S #:		
16.405	120641	14		7	0		0	95
Kexadeca	noic acid				CA	s #: 57-10-3		
16.493	61138	7	•	4	97	NBS75K.l	71609	95
hknown					CA	us #:	ı	
19.864	210340	27	•	14	0	·· •	0	114

EPA SAMPLE NO.

333-MW01-1B

Lab Name: CH2M HILL Contract: MB370

Lab Code: MGM Case No.: MB370 SAS No.: SDG No.: MB370

Matrix: (soil/water) WATER Lab Sample ID: MB370005

Sample wt/vol: 1000 (g/mL) ML Lab File ID: H0006690.D

Level: (low/med) LOW Date Received: 07/13/96

% Moisture: not dec. Date Extracted:07/16/96

Extraction: (SepF/Cont/Sonc) CONT Date Analyzed: 07/24/96

GPC Cleanup: (Y/N) N pH: 7.0 Dilution Factor: 1.0

CONCENTRATION UNITS:

CAS NO. COMPOUND (ug/L or ug/Kg) UG/L Q

62-75-9N-Nitrosodimethylamine	10	U
108-95-2Phenol	10	Ü
111-44-4bis(2-Chloroethyl)ether	10	บ
95-57-82-Chlorophenol	10	ָ ע
541-73-11,3-Dichlorobenzene	10	ם
106-46-71,4-Dichlorobenzene	10	Ü
95-50-11,2-Dichlorobenzene	10	Ü
108-60-12,2'-oxybis(1-Chloroprop_(1)	10	ט
621-64-7N-Nitroso-di-n-propylamine	10	ט
67-72-1Hexachloroethane	10	מ
98-95-3Nitrobenzene		ש
	10	
78-59-1Isophorone	10	Ŭ
88-75-52-Nitrophenol	10	ָ <u>.</u>
105-67-92,4-Dimethylphenol	10	Ŭ
111-91-1bis(2-Chloroethoxy)methane	10	Ū
120-83-22,4-Dichlorophenol	10	บ
120-82-11,2,4-Trichlorobenzene	10	ָּט
91-20-3Naphthalene	10	ַ
87-68-3Hexachlorobutadiene	10	U
59-50-74-Chloro-3-methylphenol	10	ָ ַ ע
88-06-22,4,6-Trichlorophenol	10	´ U
91-58-72-Chloronaphthalene	10	U
131-11-3Dimethylphthalate	10	U
606-20-22,6-Dinitrotoluene	10	U
208-96-8Acenaphthylene	10	Ū
83-32-9Acenaphthene	10	ט
51-28-52,4-Dinitrophenol	50	ט
100-02-74-Nitrophenol	50	U
121-14-22,4-Dinitrotoluene	10	ן ע
84-66-2Diethylphthalate	10	ט
86-73-7Fluorene	10	Ū
7005-72-34-Chlorophenyl-phenylether	10	Ū

(1) 2,2'-oxybis(1-Chloropropane) is known as bis(2-Chloroisopropyl) ether

 $_{ ext{SW846}}$   $\eta$ 

FORM I SV-1

### 1C SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

333-MW01-1B

Lab Name: CH2M HILL Contract: MB370

Lab Code: MGM Case No.: MB370 SAS No.: SDG No.: MB370

Matrix: (soil/water) WATER Lab Sample ID: MB370005

Sample wt/vol: 1000 (g/mL) ML Lab File ID: H0006690.D

Level: (low/med) LOW Date Received: 07/13/96

% Moisture: not dec.\_\_\_\_ dec.\_\_ Date Extracted:07/16/96

Extraction: (SepF/Cont/Sonc) CONT Date Analyzed: 07/24/96

GPC Cleanup: (Y/N) N pH: 7.0 Dilution Factor: 1.0

CONCENTRATION UNITS:

CAS NO. COMPOUND (ug/L or ug/Kg) UG/L Q

, <del></del>		
534-52-14,6-Dinitro-2-methylphenol 86-30-6N-Nitrosodiphenylamine(3) 122-66-71,2-Diphenylhydrazine 101-55-34-Bromophenyl-phenylether 118-74-1Hexachlorobenzene 87-86-5Pentachlorophenol 85-01-8Phenanthrene 120-12-7Anthracene	50 10 10 10 10 10 50 10	ממממממם
120-12-7Anthracene 84-74-2Di-n-butylphthalate	10	U U
206-44-0Fluoranthene	10 50	ָּט
129-00-0Pyrene	10	ט
85-68-7Butylbenzylphthalate 56-55-3Benzo(a)anthracene	10 10	บ บ
91-94-13,3'-Dichlorobenzidine 218-01-9Chrysene	20 10	ប
117-81-7bis(2-Ethylhexyl)phthalate	10 10	Ū
205-99-2Benzo (b) fluoranthene 207-08-9Benzo (k) fluoranthene	10 10	ט "ע
50-32-8Benzo (a) pyrene 193-39-5Indeno (1,2,3-cd) pyrene	10 10	ט
53-70-3Dibenz(a,h)anthracene 191-24-2Benzo(g,h,i)perylene	10 10	U U
		İ

(3) - Cannot be separated from Diphenylamine

FORM I SV-2



4 15

333-MW01-1B

Lab Name: CH2M HILL

Contract: MB370

Lab Code: MGM

Case No.: MB370 SAS No.: SDG No.: MB370

Matrix: (soil/water) WATER

Lab Sample ID: MB370005

Sample wt/vol: 1000 (g/mL) ML

Lab File ID: H0006690.D

Date Received: 07/13/96

Level: (low/med) LOW

Date Extracted:07/16/96

Extraction: (SepF/Cont/Sonc) CONT

Date Analyzed: 07/24/96

GPC Cleanup: (Y/N) N

% Moisture: not dec.\_\_\_\_ dec.

pH: 7.0

Dilution Factor: 1.0

CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L

Number TICs found: 11

		<u>.                                    </u>		
CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	Unknown	5.277	3	J
2.	Unknown	5.439	2	J
3.	Unknown Unknown	5.483 5.798	3	J
4.	Unknown	6.596	4	J
6. 90-05-1	Phenol, 2-methoxy-	7.512	4	IJ
7. 65-85-0	Benzoic Acid	8.033	3	ŲИ
8. 13679-75-9 9. 121-33-5	1-(2-Thienyl)-1-propanone Vanillin	8.524 10.898	3	NJ NJ
10.	Unknown	12.649	3	J
1	Unknown	16.555	4	J
11. 12. 13.				
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1 1.53 .				
16.				
16. 17.				
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41.				
1 44.				
23.				
1 43.				
1 20.		<u> </u>		
27. 28.				
1 43.				
30.				
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SW846  $\eta$ 

Page 1

CONCENTRATIONS

Report Date: 31-Jul-1996 14:32

## CH2M Hill Montgomery

## SW846 SEMI-VOLATILES

Data file : /chem/ms3.i/h072496a.b/h0006690.d / Lab Smp Id: MB370005 / Client Client Smp ID: 333-MW01-1B

Inj Date : 24-JUL-1996 19:59

Operator : mjohnson Inst ID: ms3.i

Smp Info : MB370005 333-MW01-1B

Misc Info : Comment

Method : /chem/ms3.i/h072496a.b/SV8270.m <

Meth Date: 31-Jul-1996 14:13 mjohnson Quant Type: ISTD Cal Date: 24-JUL-1996 08:55 Cal File: h00066 Cal File: h0006673.d 🗸

Als bottle: 19

Dil Factor: 1.000~

Integrator: HP RTE Compound Sublist: brown&root.sub ~

Target Version: 3.10

		•				CONCENTRA	4110N2
			QUANT SIG			ON-COLUMN	FINAL
	Co	mpounds	MASS	RT EXP RT RE	L RT RESPONSE	( NG)	( ug/L)
	==	**************	====	== ====================================	==== ======		======
		2 N-Nitrosodimethylamine	74.00	Compound Not	Detected.		
	\$	6 2-Fluorophenol	112.00	5.549 5.550 (0	.811) 248489	150	77
		9 Phenol-d5	99.00	6.377 6.378 (0	.931) 299882	140	73
		10 Phenol	94.00	Compound Not	Detected.		
		12 bis(2-Chloroethyl)ether	63.00	Compound Not	Detected.		
		14 2-Chlorophenol	128.00	Compound Not	Detected.		
1		15 1,3-Dichlorobenzene	146.00	Compound Not	Detected.		
	*	16 1,4-Dichlorobenzene-d4	152.00	6.846 6.839 (1	.000) 41903	40	(Q)
		17 1,4-Dichlorobenzene	146.00	Compound Not	Detected.		
		20 1,2-Dichlorobenzene	146.00	Compound Not	Detected.		
		22 2,2'-oxybis(1-Chloropropane)	45.00	Compound Not	Detected.		
		26 N-Nitroso-di-n-propylamine	70.00	Compound Not	Detected.		
		30 Hexachloroethane	117.00	Compound Not	Detected.		
	\$	31 Nitrobenzene-d5	82.00	7.564 7.565 (0	.876) 121605	87	44
		32 Nitrobenzene	77.00	Compound Not	Detected.		
		34 Isophorone	82.00	Compound Not	Detected.		
		35 2-Nitrophenol	139.00	Compound Not	Detected.		
		36 2,4-Dimethylphenol	107.00	Compound Not	Detected.		
		39 2,4-Dichlorophenol	162.00	Compound Not	Detected.		
		41 1,2,4-Trichlorobenzene	180.00	Compound Not	Detected.		
	*	42 Naphthalene-d8	136.00	8.634 8.635 (1	.000) 153885	40	
		43 Naphthalene	128.00	Compound Not	Detected.		
		47 Hexachlorobutadiene	225.00	Compound Not	Detected.		
		50 4-Chloro-3-methylphenol	107.00	Compound Not	Detected.		
		55 2,4,6-Trichlorophenol	196.00	Compound Not	Detected.		
	\$	57 2-Fluorobiphenyl	172.00	10.583 10.576 (0	.876) 202210	94	47
		59 2-Chloronaphthalene	162.00	Compound Not	Detected.		
		63 Dimethylphthalate	163.00	Compound Not	Detected.		
		65 2,6-Dinitrotoluene	165.00	Compound Not	Detected.	4	

Data File: /chem/ms3.i/h072496a.b/h0006690.d Report Date: 31-Jul-1996 14:32

						CONCENT	RATIONS
		QUANT SIG		•		ON-COLUMN	FINAL
Compo	unds	MASS	RT	EXP RT REL RT	RESPONSE	( NG)	( ug/L)
=====	=======================================	====	==		=======	======	======
66	Acenaphthylene	152.00	Con	apound Not Detect	ed.		
* 68	Acenaphthene-d10	164.00	12.078	12.079 (1.000)	85491	40	
69	Acenaphthene	154.00	Cor	mpound Not Detect	ed.		
70	2,4-Dinitrophenol	184.00	Con	mpound Not Detect	ed.		
71	4-Nitrophenol	65.00	Con	mpound Not Detect	ed.		
74	2,4-Dinitrotoluene	165.00	Con	npound Not Detect	ed.		
78	Diethylphthalate	149.00	Con	npound Not Detect	ed.		
79	Fluorene	166.00	Cor	npound Not Detect	ed.		
80	4-Chlorophenyl-phenylether	204.00	Cor	mpound Not Detect	ed.		
83	4,6-Dinitro-2-methylphenol	198.00	Cor	npound Not Detect	ed.		
84	N-Nitrosodiphenylamine (1)	169.00	Cor	mpound Not Detect	ed.		
85	1,2-Diphenylhydrazine	77.00	Cor	mpound Not Detect	ed.		
\$ 86	2,4,6-Tribromophenol	329.65	13.888	13.889 (1.150)	61348	140	72
89	4-Bromophenyl-phenylether	248.00	Cor	mpound Not Detect	ed.		
90	<b>Hexach</b> lorobenzene	284.00	Cor	mpound Not Detect	ed.		
92	Pentachlorophenol	266.00	Cor	mpound Not Detect	ed.		
* 95	Phenanthrene-d10	188.00	15.492	15.501 (1.000)	124343	40	
96	Phenanthrene	178.00	Cor	mpound Not Detect	ed.		
97	Anthracene	178.00	Cor	mpound Not Detect	ed.		
99	Di-n-butylphthalate	149.00	Cor	mpound Not Detect	ed.		
102	Fluoranthene	202.00	Cor	mpound Not Detect	ed.		
103	Benzidine	184.00	Cor	mpound Not Detect	ed.		
104	Pyrene	202.00	Cor	mpound Not Detect	ed.		
106	Terphenyl-d14	244.00	19.354	19.348 (0.880)	234555	74	37
110	Butylbenzylphthalate	149.00	Cor	mpound Not Detect	ed.		
112	Benzo(a)anthracene	228.00	Cor	mpound Not Detect	ed.		
113	3,3'-Dichlorobenzidine	252.00	Cor	mpound Not Detect	ed.		
* 114	Chrysene-d12	240.00	21.999	22.008 (1.000)	99086	40	
115	Chrysene	228.00	Cor	mpound Not Detect	ed.		
116	bis(2-Ethylhexyl)phthalate	149.00	Cor	mpound Not Detect	ed.		
117	Di-n-octylphthalate	149.00	Çor	apound Not Detect	ed.		
118	Benzo(b)fluoranthene	252.00	Cor	mpound Not Detect	ed.		
120	Benzo(k)fluoranthene	252.00	Сон	mpound Not Detect	ed.		
122	Benzo(a)pyrene	252.00	Cor	ipound Not Detect	ed.		
* 123	Perylene-d12	264.00	25.509	25.510 (1.000)	96455	40	
125	Indeno(1,2,3-cd)pyrene	276.00	Cor	mpound Not Detect	ed.		
126	Dibenz(a,h)anthracene	278.00	Cor	mpound Not Detect	ed.		
127	Benzo(g,h,i)perylene	276.00	Cor	mpound Not Detect	ed.		

# QC Flag Legend

Q - Qualifier signal failed the ratio test.

\$ 33 W

Data File: /chem/ms3.i/h072496a.b/h0006690.d

Report Date: 31-Jul-1996 14:32

Page 3

# CH2M Hill Montgomery

Unknown Compounds Quantitation Report

Data file : /chem/ms3.i/h072496a.b/h0006690.d

Lab Smp Id: MB370005 Client Smp ID: 333-MW01-1B

Inj Date : 24-JUL-1996 19:59

Inst ID: ms3.i Operator : mjohnson

Smp Info : MB370005 333-MW01-1B

Misc Info :

Comment

Method : /chem/ms3.i/h072496a.b/SV8270.m

Meth Date : 31-Jul-1996 14:13 mjohnson Cal Date : 24-JUL-1996 08:55 Cal File: h0006673.d

Als bottle: 19

Dil Factor: 1.000 Target Version: 3.10

Integrator: HP RTE Compound Sublist: brown&root.sub

Sample Matrix: WATER

Quantitative Mode : Use RF of Nearest Std

IS ==	TD	RT ====	AREA	AMOUNT
*	16 1,4-Dichlorobenzene-d4	6.846	298247	40.000
*	42 Naphthalene-d8	8.634	355512	40.000
	68 Acenaphthene-d10	12.078	385706	40.000
*	95 Phenanthrene-d10	15.492	367883	40.000

		CONCEN	ITRATIONS		Q	UANT	
RT	AREA	ON-COL( NG)	FINAL( ug/L)	QUAL	LIBRARY	LIB ENTRY	CPND #
====	====	=========	. **********	====	======	========	=====
	•	110196					
Unknown	Atcohol ()	s florate		CAS	#:		
5.277	43475	6	3	0		0	16
Unknown				CAS	#:		
5.439	34534	. 5	2	0		0	16
Unknown				CAS	#:		
5.483	40618	Robling 5	3	0		0	16
		eller.					
Unknown	A <del>lcohul</del>	Upo,		CAS	#:		
5.798	89837		6	0		0	16

		CON	ICENTRAT I	ONS			QL	JANT		
RT	AREA	ON-COL(	NG) FI	NAL( ug/L)	QUAL		LIBRARY	LIB ENTRY	CPND #	
====	2255	========	=== ==	=======	====		======	=======	=====	
Unknown-Al	cohol					CAS	#:			
6.596	56962	8	3	4	0			0	16	
Phenol, 2-	methoxy-					CAS	#: 90-05-1			
7.512	60419	8	3	4	91		NBS75K.l	64772	16	
Benzoic Ad	id .					CAS	#: 65-85-0			
8.033	51612		5	3	97		NBS75K.l	3903	42	
1-(2-Thier	nyl)-1-pr	opanone				CAS	#: 13679-75	.9		
8.524	54316	. 6	5	3	83		NBS75K.l	7304	42	
Vanillin						CAS	#: 121-33-5			
10.898	111390	. 12	2	6	98		NBS75K.l	66916	68	
Unknown						CAS	#:			
12.649	64701	7	7	3	0			0	68	
Unknown			-			CAS	#:			
16.555	74634	8	3	4	0			0	95	

333-MW03-1D

Lab Name: CH2M HILL

Contract: MB370

Lab Code: MGM Case No.: MB370 SAS No.: SDG No.: MB370

Matrix: (soil/water) WATER Lab Sample ID: MB370006

Sample wt/vol: 1000 (g/mL) ML Lab File ID: H0006703.D

Level: (low/med) LOW Date Received: 07/13/96

% Moisture: not dec.\_\_\_\_\_ dec.\_\_\_ Date Extracted:07/16/96

Extraction: (SepF/Cont/Sonc) CONT Date Analyzed: 07/25/96

GPC Cleanup: (Y/N) N pH: 7.0 Dilution Factor: 1.0

CONCENTRATION UNITS: CAS NO. COMPOUND (ug/L or ug/Kg) UG/L Q

108-95-2Phenol 111-44-4bis(2-Chloroethyl)ether 95-57-82-Chlorophenol 10541-73-11,3-Dichlorobenzene 106-46-71,4-Dichlorobenzene 1095-50-11,2-Dichlorobenzene 1008-60-12,2'-oxybis(1-Chloroprop(1)) 621-64-7Nritroso-di-n-propylamine 67-72-1Hexachloroethane 1098-95-3Isophorone 1098-95-3Isophorone 1008-67-92,4-Dimethylphenol 11091-1	<del></del>		
108-95-2Phenol 111-44-4bis (2-Chloroethyl) ether 95-57-82-Chlorophenol 10541-73-11, 3-Dichlorobenzene 106-46-71, 4-Dichlorobenzene 10 95-50-11, 2-Dichlorobenzene 10 108-60-12, 2'-oxybis (1-Chloroprop (1) 621-64-7N-Nitroso-di-n-propylamine 67-72-1Hexachloroethane 98-95-3Nitrobenzene 10 98-95-3Isophorone 10 88-75-52-Nitrophenol 105-67-92, 4-Dimethylphenol 110-91-1bis (2-Chloroethoxy) methane 120-83-22, 4-Dichlorophenol 120-82-11, 2, 4-Trichlorobenzene 10 91-20-3Naphthalene 87-68-3	62-75-9N-Nitrosodimethylamine	10	U
111-44-4			บ
95-57-82-Chlorophenol			บั
541-73-11, 3-Dichlorobenzene       10         106-46-71, 4-Dichlorobenzene       10         95-50-11, 2-Dichlorobenzene       10         108-60-12, 2'-oxybis (1-Chloroprop_(1)       10         621-64-7N-Nitroso-di-n-propylamine       10         67-72-1Hexachloroethane       10         98-95-3Nitrobenzene       10         78-59-1	95-57-82-Chlorophenol		Ü
106-46-71, 4-Dichlorobenzene       10         95-50-11, 2-Dichlorobenzene       10         108-60-12, 2'-oxybis (1-Chloroprop_(1))       10         621-64-7N-Nitroso-di-n-propylamine       10         67-72-1Hexachloroethane       10         98-95-3Nitrobenzene       10         78-59-1Isophorone       10         88-75-52-Nitrophenol       10         105-67-92, 4-Dimethylphenol       10         111-91-1bis (2-Chloroethoxy) methane       10         120-83-22, 4-Dichlorophenol       10         120-82-11, 2, 4-Trichlorobenzene       10         91-20-3Naphthalene       10         87-68-3			บ
95-50-11, 2-Dichlorobenzene       10         108-60-12, 2'-oxybis (1-Chloroprop_(1))       10         621-64-7N-Nitroso-di-n-propylamine       10         67-72-1Hexachloroethane       10         98-95-3Nitrobenzene       10         78-59-1Isophorone       10         88-75-5			Ŭ
108-60-12,2'-oxybis (1-Chloroprop_(1)       10         621-64-7Nhitroso-di-n-propylamine       10         67-72-1Hexachloroethane       10         98-95-3Nitrobenzene       10         78-59-1Isophorone       10         88-75-52-Nitrophenol       10         105-67-92,4-Dimethylphenol       10         111-91-1bis (2-Chloroethoxy) methane       10         120-83-22,4-Dichlorophenol       10         120-82-11,2,4-Trichlorobenzene       10         91-20-3Naphthalene       10         87-68-3Naphthalene       10         87-68-3			Ū
621-64-7Nitroso-di-n-propylamine       10         67-72-1Hexachloroethane       10         98-95-3Nitrobenzene       10         78-59-1Isophorone       10         88-75-5			ับ
67-72-1			บ
98-95-3Nitrobenzene       10         78-59-1Isophorone       10         88-75-52-Nitrophenol       10         105-67-92,4-Dimethylphenol       10         111-91-1bis(2-Chloroethoxy)methane       10         120-83-22,4-Dichlorophenol       10         120-82-11,2,4-Trichlorobenzene       10         91-20-3Naphthalene       10         87-68-3Hexachlorobutadiene       10         59-50-74-Chloro-3-methylphenol       10         88-06-22,4,6-Trichlorophenol       10         91-58-72-Chloronaphthalene       10         131-11-3Dimethylphthalate       10         606-20-22,6-Dinitrotoluene       10         208-96-8Acenaphthylene       10         83-32-9Acenaphthene       50         100-02-74-Nitrophenol       50         121-14-22,4-Dinitrotoluene       10		ı	Ū
78-59-1			บี
88-75-52-Nitrophenol       10         105-67-92,4-Dimethylphenol       10         111-91-1bis (2-Chloroethoxy) methane       10         120-83-22,4-Dichlorophenol       10         120-82-11,2,4-Trichlorobenzene       10         91-20-3Naphthalene       10         87-68-3Hexachlorobutadiene       10         59-50-74-Chloro-3-methylphenol       10         88-06-22,4,6-Trichlorophenol       10         91-58-72-Chloronaphthalene       10         131-11-3Dimethylphthalate       10         606-20-22,6-Dinitrotoluene       10         208-96-8Acenaphthene       10         51-28-52,4-Dinitrophenol       50         100-02-74-Nitrophenol       50         121-14-22,4-Dinitrotoluene       10			Ü
105-67-92,4-Dimethylphenol       10         111-91-1bis (2-Chloroethoxy) methane       10         120-83-22,4-Dichlorophenol       10         120-82-11,2,4-Trichlorobenzene       10         91-20-3Naphthalene       10         87-68-3Hexachlorobutadiene       10         59-50-74-Chloro-3-methylphenol       10         88-06-22,4,6-Trichlorophenol       10         91-58-72-Chloronaphthalene       10         131-11-3Dimethylphthalate       10         606-20-22,6-Dinitrotoluene       10         208-96-8Acenaphthylene       10         83-32-9Acenaphthene       10         51-28-52,4-Dinitrophenol       50         100-02-74-Nitrophenol       50         121-14-22,4-Dinitrotoluene       10			Ü
111-91-1			บี
120-83-22,4-Dichlorophenol       10         120-82-11,2,4-Trichlorobenzene       10         91-20-3Naphthalene       10         87-68-3Hexachlorobutadiene       10         59-50-74-Chloro-3-methylphenol       10         88-06-22,4,6-Trichlorophenol       10         91-58-72-Chloronaphthalene       10         131-11-3Dimethylphthalate       10         606-20-22,6-Dinitrotoluene       10         208-96-8Acenaphthylene       10         83-32-9Acenaphthene       10         51-28-52,4-Dinitrophenol       50         100-02-74-Nitrophenol       50         121-14-22,4-Dinitrotoluene       10			บั
120-82-11, 2, 4-Trichlorobenzene       10         91-20-3Naphthalene       10         87-68-3Hexachlorobutadiene       10         59-50-74-Chloro-3-methylphenol       10         88-06-22, 4, 6-Trichlorophenol       10         91-58-72-Chloronaphthalene       10         131-11-3Dimethylphthalate       10         606-20-22, 6-Dinitrotoluene       10         208-96-8Acenaphthylene       10         83-32-9Acenaphthene       10         51-28-52, 4-Dinitrophenol       50         100-02-74-Nitrophenol       50         121-14-22, 4-Dinitrotoluene       10			Ŭ
91-20-3Naphthalene       10         87-68-3Hexachlorobutadiene       10         59-50-74-Chloro-3-methylphenol       10         88-06-22,4,6-Trichlorophenol       10         91-58-72-Chloronaphthalene       10         131-11-3Dimethylphthalate       10         606-20-22,6-Dinitrotoluene       10         208-96-8Acenaphthylene       10         83-32-9Acenaphthene       10         51-28-52,4-Dinitrophenol       50         100-02-74-Nitrophenol       50         121-14-22,4-Dinitrotoluene       10			บี
87-68-3			Ŭ
59-50-74-Chloro-3-methylphenol       10         88-06-22,4,6-Trichlorophenol       10         91-58-72-Chloronaphthalene       10         131-11-3Dimethylphthalate       10         606-20-22,6-Dinitrotoluene       10         208-96-8Acenaphthylene       10         83-32-9Acenaphthene       10         51-28-52,4-Dinitrophenol       50         100-02-74-Nitrophenol       50         121-14-22,4-Dinitrotoluene       10			บี
88-06-22,4,6-Trichlorophenol       10         91-58-72-Chloronaphthalene       10         131-11-3Dimethylphthalate       10         606-20-22,6-Dinitrotoluene       10         208-96-8Acenaphthylene       10         83-32-9Acenaphthene       10         51-28-52,4-Dinitrophenol       50         100-02-74-Nitrophenol       50         121-14-22,4-Dinitrotoluene       10			บี
91-58-72-Chloronaphthalene       10         131-11-3Dimethylphthalate       10         606-20-22,6-Dinitrotoluene       10         208-96-8Acenaphthylene       10         83-32-9Acenaphthene       10         51-28-52,4-Dinitrophenol       50         100-02-74-Nitrophenol       50         121-14-22,4-Dinitrotoluene       10			
131-11-3Dimethylphthalate       10         606-20-22,6-Dinitrotoluene       10         208-96-8Acenaphthylene       10         83-32-9Acenaphthene       10         51-28-52,4-Dinitrophenol       50         100-02-74-Nitrophenol       50         121-14-22,4-Dinitrotoluene       10			Ü
606-20-22,6-Dinitrotoluene       10         208-96-8Acenaphthylene       10         83-32-9Acenaphthene       10         51-28-52,4-Dinitrophenol       50         100-02-74-Nitrophenol       50         121-14-22,4-Dinitrotoluene       10			Ŭ
208-96-8			Ŭ
83-32-9Acenaphthene       10         51-28-52,4-Dinitrophenol       50         100-02-74-Nitrophenol       50         121-14-22,4-Dinitrotoluene       10			Ŭ
51-28-52,4-Dinitrophenol       50         100-02-74-Nitrophenol       50         121-14-22,4-Dinitrotoluene       10			บี
100-02-74-Nitrophenol 50 121-14-22,4-Dinitrotoluene 10			Ŭ
121-14-22,4-Dinitrotoluene10			Ŭ
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SW846 mm (1) 2,2'-oxybis(1-Chloropropane) is known as bis(2-Chloroisopropyl) ether

EPA SAMPLE NO.

333-MW03-1D

Lab Name: CH2M HILL Contract: MB370

Lab Code: MGM Case No.: MB370 SAS No.: SDG No.: MB370

Matrix: (soil/water) WATER Lab Sample ID: MB370006

Sample wt/vol: 1000 (g/mL) ML Lab File ID: H0006703.D

Level: (low/med) LOW Date Received: 07/13/96

% Moisture: not dec.\_\_\_\_\_ dec.\_\_\_\_ Date Extracted:07/16/96

Extraction: (SepF/Cont/Sonc) CONT Date Analyzed: 07/25/96

GPC Cleanup: (Y/N) N pH: 7.0 Dilution Factor: 1.0

CONCENTRATION UNITS:

CAS NO. COMPOUND (ug/L or ug/Kg) UG/L Q

	<del></del>	
534-52-14,6-Dinitro-2-methylphenol	50	U
86-30-6N-Nitrosodiphenylamine (3)	10	U
122-66-71,2-Diphenylhydrazine	10	U
101-55-34-Bromophenyl-phenylether	10	ט
118-74-1Hexachlorobenzene	10	ן ט
87-86-5Pentachlorophenol	50	U
85-01-8Phenanthrene	10	ַ ע
120-12-7Anthracene	10	U
84-74-2Di-n-butylphthalate	10	U
206-44-0Fluoranthene	10	U
92-87-5Benzidine	50	U
129-00-0Pyrene	10	U
85-68-7Butylbenzylphthalate	10	U
56-55-3Benzo (a) anthracene	10	U
91-94-13,3'-Dichlorobenzidine	20	U
218-01-9Chrysene	10	U
117-81-7bis(2-Ethylhexyl)phthalate	10	Ŭ
117-84-0Di-n-octylphthalate	10	Ŭ
205-99-2Benzo (b) fluoranthene	10	Ū
207-08-9Benzo(k)fluoranthene	10	្វ
50-32-8Benzo(a)pyrene	10	U
193-39-5Indeno(1,2,3-cd)pyrene	10	Ū
53-70-3Dibenz(a,h)anthracene	10	_
191-24-2Benzo(g,h,i)perylene	10	บ
	l	

(3) - Cannot be separated from Diphenylamine

FORM I SV-2



4 11

## 1F SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

333-MW03-1D

Lab Name: CH2M HILL

Contract: MB370

Case No.: MB370 SAS No.: SDG No.: MB370

Matrix: (soil/water) WATER

Lab Sample ID: MB370006

Sample wt/vol:

Lab Code: MGM

1000 (g/mL) ML

Lab File ID: H0006703.D

Level: (low/med) LOW

Date Received: 07/13/96

% Moisture: not dec. dec.

Date Extracted:07/16/96

Extraction: (SepF/Cont/Sonc) CONT

Date Analyzed: 07/25/96

GPC Cleanup: (Y/N) N

pH: 7.0

Dilution Factor: 1.0

CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L

Number TICs found: 7

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1. 123-42-2 2. 111-76-2 3. 111-90-0 4. 13429-07-7 5. 6. 7. 8. 9. 10. 11. 12. 13. 14. 15. 16. 17. 18. 19. 20. 21. 22. 23. 24. 25. 26. 27. 28. 29.	2-Pentanone, 4-hydroxy-4-met Ethanol, 2-butoxy- Ethanol, 2-(2-ethoxyethoxy)- 2-Propanol, 1-(2-methoxyprop Unknown Hydrocarbon Unknown Unknown	5.182 5.717 6.516		Q NJAB NJB NJ J J J J J J J J J J J J J J J J
28. 29. 30.				

Report Date: 05-Aug-1996 14:11

## CH2M Hill Montgomery

SW846 SEMI-VOLATILES

Data file : /chem/ms3.i/h072596a.b/h0006703.d/ Lab Smp Id: MB370006 Client Client Smp ID: 333-MW03-1D

Inj Date : 25-JUL-1996 16:30

Operator : mjohnson Inst ID: ms3.i.

Smp Info : MB370006 333-MW03-1D

Misc Info : Comment

Method : /chem/ms3.i/h072596a.b/SV8270.m /

Meth Date : 05-Aug-1996 14:10 mjohnson Quant Type: ISTD Cal Date : 25-JUL-1996 09:31 Cal File: h000669 Cal File: h0006693.d —

Als bottle: 12

Dil Factor: 1.000 ~

Integrator: HP RTE Compound Sublist: brown&root.sub /

Target Version: 3.10

		_					CONCENTR	ATIONS
		•	QUANT SIG				ON-COLUMN	FINAL
Co	mpounds		MASS	RT	EXP RT REL RT	RESPONSE	( NG)	( ug/L)
==		====	====	==		=======	======	======
	2 N-Nitrosodimeth	ylamine	74.00	Com	pound Not Detect	ed.		
\$	6 2-Fluorophenol		112.00	5.468	5.471 (0.810)	264332	160	82
\$	9 Phenol-d5		99.00	6.289	6.291 (0.932)	315334	160	78
	:10 Phenol		94.00	Com	pound Not Detect	ed.		
	12 bis(2-Chloroeth	yl)ether	63.00	Com	pound Not Detect	ed.		
	14 2-Chlorophenol		128.00	Com	pound Not Detect	ed.		
	15 1,3-Dichloroben	zene	146.00	Com	pound Not Detect	ed.		
1*	16 1,4-Bichloroben	zene-d4	152.00	6.750	6.753 (1.000)	40637	40	(Q)
	17 1,4-Dichloroben	zene	146.00	Com	pound Not Detect	ed.		
	20 1,2-Dichloroben	zene	146.00	Com	pound Not Detect	ed.		
	22 2,2'-oxybis(1-C	hloropropane)	45.00	Com	pound Not Detect	ed.		
	26 N-Nitroso-di-n-	propylamine	70.00	Com	pound Not Detect	ed.		
	30 Hexachloroethan	e	117.00	Сот	pound Not Detect	ed.		
\$	31 Nitrobenzene-d5		82.00	7.461	7.464 (0.876)	128547	88	44
	32 Nitrobenzene		77.00	Com	pound Not Detect	ed.		
	34 Isophorone		82.00	Com	pound Not Detect	ed.		
	35 2-Nitrophenol		139.00	Com	pound Not Detect	ed.		
	36 2,4-Dimethylphe	nol	107.00	Соп	pound Not Detect	ed.		
	38 bis(2-Chloroeth	oxy)methane	93.00	Com	pound Not Detect	ed.		
	39 2,4-Dichlorophe	nol	162.00	Com	pound Not Detect	ed.		
,	41 1,2,4-Trichloro	benzene	180.00	Соп	pound Not Detect	ed.		
*	42 Naphthalene-d8		136.00	8.516	8.519 (1.000)	147862	40	
	43 Naphthalene		128.00	Соп	pound Not Detect	ed.		
	47 Hexachlorobutad	i ene	225.00	Com	pound Not Detect	ed.		
	50 4-Chloro-3-meth	ylphenol	107.00	Com	pound Not Detect	ed.		
	55 2,4,6-Trichloro	phenol	196.00	Com	pound Not Detect	ed.		
\$	57 2-fluorobipheny	ι	172.00	10.451	10.453 (0.875)	197936	88	44
	59 2-Chloronaphtha	l ene	162.00	Con	pound Not Detect	ed.		
	63 Dimethylphthala	te	163.00	Соп	pound Not Detect	ed.		
								۲, ۲

							CONC	ENTRATIONS
			QUANT SIG		•		ON-COL	UMN FINAL
C	ompounds		MASS	RT	EXP RT REL RT	RESPONSE	( NO	G) (ug/L)
=:		=======	====	==		=======	=====	== ======
	65 2,6-Dinitro	toluene	165.00	Cor	mpound Not Detect	ed.		
	66 Acenaphthyl	ene	152.00	Cor	mpound Not Detect	ed.		
*	68 Acenaphthen	e-d10	164.00	11.938	11.941 (1.000)	80316		40
	69 Acenaphthen	e	154.00	Co	mpound Not Detect	ed.		
	70 2,4-Dinitro	phenol	184.00	Co	mpound Not Detect	ed.		
	71 4-Nitrophen	ol	65.00	Co	mpound Not Detect	ed.		
	74 2,4-Dinitro	toluene	165.00	Со	mpound Not Detect	ed.		
	78 Diethylphth	alate	149.00	Co	mpound Not Detect	ed.		
	79 Fluorene		166.00	Co	mpound Not Detect	ed.		
	80 4-Chlorophe	nyl-phenylether	204.00	Co	mpound Not Detect	ed.		
	83 4,6-Dinitro	-2-methylphenol	198.00	Co	mpound Not Detect	ed.		
	84 N-Nitrosodi	phenylamine (1)	169.00	Co	mpound Not Detect	ed.		
	85 1,2-Dipheny	lhydrazine	77.00	Co	mpound Not Detect	ed.		
\$	86 2,4,6-Tribr	omophenol	329.65	13.741	13.744 (1.151)	59616	19	50 76
	89 4-Bromophen	yl-phenylether	248.00	Co	mpound Not Detect	ed.		
	90 Hexachlorob	enzene	284.00	Co	mpound Not Detect	ed.		
	92 Pentachloro	phenol	266.00	Co	mpound Not Detect	ed.		
*	95 Phenanthren	e-d10 .	188.00	15.339	15.348 (1.000)	111038		40
	96 Phenanthren	e	178.00	Co	mpound Not Detect	ed.		
	97 Anthracene		178.00	Co	mpound Not Detect	ed.		
	99 Di-n-butylp	hthalate	149.00	Co	mpound Not Detect	ed.		
	102 Fluoranthen	e	202.00	Co	mpound Not Detect	ed.		
	103 Benzidine		184.00	Co	mpound Not Detect	ed.		
	104 Pyrene		202.00	Co	mpound Not Detect	ed.		
÷	106 Terphenyl-d	14	244.00		19.203 (0.879)	193778	•	77 38
	110 Butylbenzyl	phthalate	149.00	Co	mpound Not Detect	ed.		
	112 Benzo(a)ant	hracene	228.00	Co	mpound Not Detect	ed.		
1	113 3,3'-Dichlo	robenzidine	252.00	Co	mpound Not Detect	ed.		
. *	114 Chrysene-d1	2	240,00		21.848 (1.000)	85881		40 -
	115 Chrysene		228.00	Co	mpound Not Detect	ted.		
	116 bis(2-Ethyl	hexyl)phthalate	149.00	Co	mpound Not Detect	ted.		
	117 Di-n-octylp	hthalate	149.00		mpound Not Detect			
	118 Benzo(b)flu		252.00		mpound Not Detect			
	120 Benzo(k)flu	oranthene	252.00		mpound Not Detect			
	122 Benzo(a)pyr	ene .	252.00		mpound Not Detect			
*	123 Perylene-d1		264.00		25.329 (1.000)	83427		40
	125 Indeno(1,2,		276.00		mpound Not Detect			- <del>-</del>
	126 Dibenz(a,h)		278.00		mpound Not Detect			
	127 Benzo(g,h,i		276.00		mpound Not Detect			
	=- 10,	•		30		<del></del>		

QC Flag Legend

Q - Qualifier signal failed the ratio test.

Data File: /chem/ms3.i/h072596a.b/h0006703.d

Report Date: 31-Jul-1996 16:15

Page 3

# CH2M Hill Montgomery

Unknown Compounds Quantitation Report

Data file: /chem/ms3.i/h072596a.b/h0006703.d

Lab Smp Id: MB370006 Client Smp ID: 333-MW03-1D

Inj Date : 25-JUL-1996 16:30

Operator : mjohnson Smp Info : MB370006 333-MW03-1D Inst ID: ms3.i

Misc Info : Comment

Method : /chem/ms3.i/h072596a.b/SV8270.m

Meth Date: 31-Jul-1996 16:11 mjohnson

Cal Date : 25-JUL-1996 09:31 Cal File: h0006693.d

Als bottle: 12

Target Version: 3.10

Dil Factor: 1.000 Integrator: HP RTE Compound Sublist: brown&root.sub

Sample Matrix: WATER

Quantitative Mode : Use RF of Nearest Std

ISTD	RT ====	AREA	AMOUNT
* 16 1,4-Dichlorobenzene-d4	6.750	263729	40.000
* 95 Phenanthrene-d10	15.339	339363	40.000
* 114 Chrysene-d12	21.838	328230	40.000

,	~	CONCEN	TRATIONS		Qt	JANT	
· RT	AREA	ON-COL( NG)	FINAL( ug/L)	QUAL	LIBRARY	LIB ENTRY	CPND #
222	====	**********	========	====	======	=======	=====
✓ 2-Pent	anone, 4-hyd	droxy-4-methyl-		CAS #	<b>#:</b> 123-42-2		
5.18	2 `50512	2 8	4	42	NBS75K.l	64274	16
Unknow	n-Atcohot			CAS #	4: 111-76	2	
5.71	7 6068	7 9	5	0		0	16
/Unknow	n Atcohel			CAS #	#: 111-90+	0	
6.51	6 84596	6 13	6	0		0	16
Unknow	n Alcohol			CAS i	#: 13 <b>+</b> 2-9	·07-7	
6.59	7 3482	3 5	3	0		0	16
Unknow	n hydroc	an lean		CAS :	<b>#:</b>		
16.32	0 5 <i>77</i> 89	9 7	3	0		0	95

Data File: /chem/ms3.i/h072596a.b/h0006703.d Report Date: 31-Jul-1996 16:15

		CONCENT	RATIONS		UANT		
RT	AREA ON-	COL( NG)	FINAL( ug/L)	QUAL	LIBRARY	LIB ENTRY	CPND #
====	2222 <b>2</b> 22	========	3========	====	222222	2322222	=====
Unknown	hydrocar	bon		CAS	· #:		
16.401	44183	5	3	0		0	95
Unknown				CAS	<b>:</b> #:		
19.852	230415	28	14	0		0	114

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GC PURGEABLE HALOCARBONS

# CASE NARRATIVE GC PURGEABLE HALOCARBONS

QAL Lab	Reference	No	./SDG	MB	370			
							•	
Project:	BROW	3 V	ROOT	COASTAL	SYSTEMS	STATION		

#### I. RECEIPT

No exceptions were encountered unless a Sample Receipt Exception Report is attached to the Chain-of-Custody included with this data package.

### II. HOLDING TIMES

- A. Sample Preparation: Not applicable.
- B. Sample Analysis: All holding times were met.

#### III. METHOD

Preparation: N/A Cleanup: N/A

Analysis: EPA 601 (MOD)

#### IV. PREPARATION

Not applicable.

#### V. ANALYSIS

- A. Calibration: All acceptance criteria were met.
- B. Blanks: All acceptance criteria were met.
- C. Surrogates: All acceptance criteria were met.
- D. Matrix Spikes: 2-Chloroethyl vinyl ether was outside acceptable limits for Accuracy (% Recovery) and Precision (RPD). However, analysis of a Laboratory Control Sample immediately after the matrix spikes indicated the analytical system was in control for the compounds found in the associated samples. Since MS/MSD results are subject to matrix effects, these values should be considered to be advisory.
- E. Samples: Sample MB370001 (333-MW01-1) contained unidentified, non-target compounds.

Primary analysis utilized a Restek Rtx 502.2 (105 meter x 0.53 mm) column. Confirmation analysis used a J&W Scientific DB-VRX (75 meter x 0.45 mm) column.

I certify that this data package is in compliance with the terms and conditions agreed to by the client and QAL, Inc., both technically and for completeness except for the conditions noted above. Release of the data contained in this hardcopy data package has been authorized by the Laboratory Manager or designated person, as verified by the following signature.

SIGNED: Well W. Smiles	DATE:	
SIGNED: Herb Kelly Organic Division Manager		

# CASE NARRATIVE Addendum

Sample Information

LAB SAMPLE ID	CLIENT SAMPLE ID	SAMPLE MATRIX	DATE SAMPLED	DATE EXTRACTED	DATE ANALYZED	SAMPLE pH <sup>1</sup>
MB370001	333-MW01-1	WATER	7/11/96	N/A	7/19/96	<2
MB370002RE	333-MW02-1	WATER	7/11/96	N/A	7/19/96	<2
MB370003	333-MW03-1	WATER	7/11/96	N/A	7/18/96	<2
MB370003MS	333-MW03-1MS	WATER	7/11/96	N/A	7/19/96	<2
MB370003MSD	333-MW03-1MD	WATER	7/11/96	N/A	7/19/96	<2
MB370004RE	333-PC4-1	WATER	7/11/96	N/A	7/19/96	<2
MB370005	333-MW01-1B	WATER	7/11/96	N/A	7/19/96	<2
MB370006	333-MW03-1D	WATER	7/11/96	N/A	7/18/96	<2
MB370007	TRIP BLANK	WATER	7/11/96	N/A	7/18/96	<2
WMV096G182	VBLK001	WATER	N/A	N/A	7/18/96	N/A
WMV096G191	VBLK002	WATER	N/A	N/A	7/19/96	N/A

Applies to samples designated for purgeable VOA analysis only.

# CURRENT METHOD DETECTION LIMITS (MDLs) PURGEABLE HALOCARBONS

Date collected: N/A Sample Group: LABQC

Date extracted: N/A Lab Sample ID: Multiple Samples

 Date analyzed:
 3/13/96
 Lab file 1 ID:
 N/A

 Matrix:
 Water
 Lab file 2 ID:
 N/A

 Method:
 EPA601 (MOD)
 Dilution factor:
 1.0

 % Moisture:
 100
 Reporting units:
 ug/L

CAS NUMBER REPORTING LIMIT RESULT COMPOUND NAME 75-27-4 Bromodichloromethane 1.0 0.093 75-25-2 Bromoform 1.0 0.142 74-83-9 Bromomethane 1.0 0.089 Carbon tetrachloride 0.090 56-23-5 1.0 108-90-7 Chlorobenzene 1.0 0.141 75-00-3 Chloroethane 1.0 0.101 110-75-8 2-Chloroethyl vinyl ether 1.0 0.100 67-66-3 Chloroform 1.0 0.086 74-87-3 Chloromethane 1.0 0.138 0.106 124-48-1 Dibromochloromethane 1.0 95-50-1 0.128 1.2-Dichlorobenzene 1.0 541-73-1 1,3-Dichlorobenzene 0.137 0.145 106-46-7 1,4-Dichlorobenzene 1.0 75-71-8 Dichlorodifluoromethane 1.0 0.181 75-34-3 1.1-Dichloroethane 1.0 0.079 107-06-2 1,2-Dichloroethane 1.0 0.089 75-35-4 1,1-Dichloroethene 1.0 0.138 0.074 156-59-2 cis-1,2-Dichloroethene 1.0 156-60-5 trans-1,2-Dichloroethene 0.066 1.0 78-87-5 1,2-Dichloropropane 1.0 0.097 10061-01-5 cis-1,3-Dichloropropene 1.0 0.095 10061-02-6 trans-1,3-Dichloropropene 1.0 0.113 75-09-2 5.0 Methylene chloride (Dichloromethane) 2.029 79-34-5 1,1,2,2-Tetrachlorethane 1.0 0.214 127-18-4 Tetrachloroethene 1.0 0.104 71-55-6 1,1,1-Trichloroethane 1.0 0.088 79-00-5 1,1,2-Trichloroethane 1.0 0.135

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1.0

1.0

1.0

79-01-6

75-69-4

75-01-4

Trichloroethene

Vinyl chloride

Trichlorofluoromethane

0.091 0.096

0.160

333-MW01-1

#### REPORT OF ANALYTICAL RESULTS PURGEABLE HALOCARBONS

Date collected: 7/11/96 Date extracted: N/A

% Moisture: 100

Date analyzed: 7/19/96 Matrix: Water Method: EPA601 (MOD)

Sample Group: MB370 Lab Sample ID: MB370001 Lab file 1 ID: G18T019 Lab file 2 ID: G18U019

Dilution factor: 1.0 Reporting units: ug/L

CAS NUMBER	COMPOUND NAME	REPORTING LIMIT	RESULT
75-27-4	Bromodichloromethane	1.0	U
75-25-2	Bromoform	1.0	ប
74-83-9	Bromomethane	1.0	U
56-23-5	Carbon tetrachloride	1.0	U
108-90-7	Chlorobenzene	1.0	υ
75-00-3 ·	Chloroethane	1.0	บ
110-75-8	2-Chloroethyl vinyl ether	1.0	ប
67-66-3	Chloroform	1.0	U
74-87-3	Chloromethane	1.0	U
124-48-1	Dibromochloromethane	1.0	U
95-50-1	1,2-Dichlorobenzene	1.0	บ
541-73-1	1,3-Dichlorobenzene	1.0	ប
106-46-7	1,4-Dichlorobenzene	1.0	ប
75-71-8	Dichlorodifluoromethane	1.0	ប
75-34-3	1,1-Dichloroethane	1.0	U
107-06-2	1,2-Dichloroethane	1.0	U
75-35-4	1,1-Dichloroethene	1.0	U
156-59-2	cis-1,2-Dichloroethene	1.0	1.8
156-60-5	trans-1,2-Dichloroethene	1.0	U
78-87-5	1,2-Dichloropropane	1.0	U
10061-01-5	cis-1,3-Dichloropropene	1.0	U
10061-02-6	trans-1,3-Dichloropropene	1.0	U
75-09-2	Methylene chloride (Dichloromethane)	5.0	. <b>U</b>
79-34-5	1,1,2,2-Tetrachlorethane	1.0	U
127-18-4	Tetrachloroethene	1.0	บ
71-55-6	1,1,1-Trichloroethane	1.0	U
79-00-5	1.1.2-Trichloroethane	1.0	U
79-01-6	Trichloroethene	1.0	U
75-69-4	Trichlorofluoromethane	1.0	U
-75-01-4	Vinyl chloride	1.0	J .
	SURROGATE-Fluorobenzene (QC Limits - 61-13:	3%)	101 % Rec.

333-MW02-1

## REPORT OF ANALYTICAL RESULTS PURGEABLE HALOCARBONS

Date collected: 7/11/96
Date extracted: N/A
Date analyzed: 7/19/96
Matrix: Water
Method: EPA601 (MOD)
% Moisture: 100

Sample Group: MB370
Lab Sample ID: MB370002RE
Lab file 1 ID: G19T017
Lab file 2 ID: G19U017
Dilution factor: 1.0

Reporting units: ug/L

CAS NUMBER	COMPOUND NAME	REPORTING LIMIT	RESULT
75-27-4	Bromodichloromethane	1.0	τ
75-25-2	Bromoform	1.0	τ
74-83-9	Bromomethane	1.0	Ţ
56-23-5	Carbon tetrachloride	1.0	τ
108-90-7	Chlorobenzene	1.0	τ
75-00-3	Chloroethane	1.0	τ
110-75-8	2-Chloroethyl vinyl ether	1.0	τ
67-66-3	Chloroform	1.0	τ
74-87-3	Chloromethane	1.0	τ
124-48-1	Dibromochloromethane	1.0	τ
95-50-1	1.2-Dichlorobenzene	1.0	τ
541-73-1 -	1.3-Dichlorobenzene	1.0	Ţ
106-46-7	1.4-Dichlorobenzene	1.0	τ
75-71-8	Dichlorodifluoromethane	1.0	Ţ
75-34-3	1.1-Dichloroethane	1.0	į
107-06-2	1,2-Dichloroethane	1.0	t
75-35-4	1,1-Dichloroethene	1.0	t
156-59-2	cis-1,2-Dichloroethene	1.0	Į
156-60-5	trans-1,2-Dichloroethene	1.0	τ
78-87-5	1,2-Dichloropropane	1.0	τ
10061-01-5	cis-1,3-Dichloropropene	1.0	τ
10061-02-6	trans-1,3-Dichloropropene	1.0	Ţ
75-09-2	Methylene chloride (Dichloromethane)	5.0	τ
79-34-5	1,1,2,2-Tetrachlorethane	1.0	τ
127-18-4	Tetrachloroethene	1.0	Ţ
71-55-6	1,1,1-Trichloroethane	1.0	Ţ
79-00-5	1,1,2-Trichloroethane	1.0	t
79-01-6	Trichloroethene	1.0	τ
75-69-4	Trichlorofluoromethane	1.0	Ţ
75-01-4	Vinyl chloride	1.0	٦ .
	SURROGATE-Fluorobenzene (QC Limits - 61-133	3%)	99 % Rec



333-MW03-1

## REPORT OF ANALYTICAL RESULTS PURGEABLE HALOCARBONS

Date collected: 7/11/96
Date extracted: N/A
Date analyzed: 7/18/96

Matrix: Water

Method: EPA601 (MOD) % Moisture: 100

Sample Group: MB370 Lab Sample ID: MB370003

Lab file 1 ID: G18T016
Lab file 2 ID: G18U016

Dilution factor: 1.0 Reporting units: ug/L

CAS NUMBER	COMPOUND NAME	REPORTING LIMIT	RESULI
75-27-4	Bromodichloromethane	1.0	τ
75-25-2	Bromoform	1.0	τ
74-83-9	Bromomethane	1.0	τ
56-23-5	Carbon tetrachloride	1.0	τ
108-90-7	Chlorobenzene	1.0	τ
75-00-3	Chloroethane	1.0	ŧ
110-75-8	2-Chloroethyl vinyl ether	1.0	τ
67-66-3	Chloroform	1.0	τ
74-87-3	Chloromethane	1.0	τ
124-48-1	Dibromochloromethane	1.0	τ
95-50-1	1,2-Dichlorobenzene	1.0	τ
541-73-1	1,3-Dichlorobenzene	1.0	τ
106-46-7	1,4-Dichlorobenzene	1.0	
75-71-8	Dichlorodifluoromethane	1.0	Ţ
75-34-3	1,1-Dichloroethane	1.0	· 1
107-06-2	1,2-Dichloroethane	1.0	τ
75-35-4	1,1-Dichloroethene	1.0	τ
156-59-2	cis-1,2-Dichloroethene	1.0	τ
156-60-5	trans-1,2-Dichloroethene	1.0	τ
78-87-5	1,2-Dichloropropane	1.0	τ
10061-01-5	cis-1,3-Dichloropropene	1.0	Ţ
10061-02-6	trans-1,3-Dichloropropene	1.0	τ
75-09-2	Methylene chloride (Dichloromethane)	5.0	t
79-34-5	1,1,2,2-Tetrachlorethane	1.0	Ţ
127-18-4	Tetrachloroethene	1.0	Ţ
71-55-6	1,1,1-Trichloroethane	1.0	τ
79-00-5	1.1.2-Trichloroethane	1.0	τ
79-01-6	Trichloroethene	1.0	τ
75-69-4	Trichlorofluoromethane	1.0	1
75-01-4	Vinyl chloride	1.0	
	SURROGATE-Fluorobenzene (QC Limits - 61-13	33%)	99 % Rec

ml

333-MW03-1MS

#### REPORT OF ANALYTICAL RESULTS PURGEABLE HALOCARBONS

Date collected: 7/11/96 Date extracted: N/A Date analyzed: 7/19/96

% Moisture: 100

Matrix: Water Method: EPA601 (MOD) Sample Group: MB370

Lab Sample ID: MB370003MS Lab file 1 ID: G18T026 Lab file 2 ID: G18U026

Dilution factor: 1.0 Reporting units: ug/L

CAS NUMBER	COMPOUND NAME F	REPORTING LIMIT	RESULT
75-27-4	Bromodichloromethane	1.0	20
75-25-2	Bromoform	1.0	20
74-83-9	Bromomethane	1.0	20
56-23-5	Carbon tetrachloride	1.0	19
108-90-7	Chlorobenzene	1.0	20
75-00-3	Chloroethane	1.0	20
110-75-8	2-Chloroethyl vinyl ether	1.0	U
67-66-3	Chloroform	1.0	20
74-87-3	Chloromethane	1.0	21
124-48-1	Dibromochloromethane	1.0	21
95-50-1	1,2-Dichlorobenzene	1.0	20
541-73-1	1,3-Dichlorobenzene	1.0	20
106-46-7	1,4-Dichlorobenzene	1.0	19
75-71-8	Dichlorodifluoromethane	1.0	20
75-34-3	1,1-Dichloroethane	1.0	20
107-06-2	1,2-Dichloroethane	1.0	20
75-35-4	1,1-Dichloroethene	1.0	19
156-59-2	cis-1,2-Dichloroethene	1.0	20
156-60-5	trans-1,2-Dichloroethene	1.0	20
78-87-5	1,2-Dichloropropane	1.0	20
10061-01-5	cis-1,3-Dichloropropene	1.0	20
10061-02-6	trans-1,3-Dichloropropene	1.0	20
75-09-2	Methylene chloride (Dichloromethane)	5.0	19
79-34-5	1,1,2,2-Tetrachlorethane	1.0	20
127-18-4	Tetrachloroethene	1.0	20
71-55-6	1,1,1-Trichloroethane	1.0	20
79-00-5	1,1,2-Trichloroethane	1.0	20
79-01-6	Trichloroethene	1.0	20
75-69-4	Trichlorofluoromethane	1.0	20
75-01-4	Vinyl chloride	1.0	- 20
	SURROGATE-Fluorobenzene (QC Limits - 61-133	3%)	100 % Rec.

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333-MW03-1MD

## REPORT OF ANALYTICAL RESULTS PURGEABLE HALOCARBONS

Date collected: 7/11/96 Sample Group: MB370

Date extracted: N/A Lab Sample ID: MB370003MSD
Date analyzed: 7/19/96 Lab file 1 ID: G18T027
Matrix: Water Lab file 2 ID: G18U027

Matrix: Water Lab file 2 ID: G18U02

Method: EPA601 (MOD) Dilution factor: 1.0

% Moisture: 100 Reporting units: ug/L

CAS NUMBER	COMPOUND NAME R	REPORTING LIMIT	RESULT
75-27-4	Bromodichloromethane	1.0	20
75-25-2	Bromoform	1.0	21
74-83-9	Bromomethane	1.0	19
56-23-5	Carbon tetrachloride	1.0	19
108-90-7	Chlorobenzene	1.0	21
75-00-3	Chloroethane	1.0	20
110-75-8	2-Chloroethyl vinyl ether	1.0	U
67-66-3	Chloroform	1.0	20
74-87-3	Chloromethane	1.0	20
124-48-1	Dibromochloromethane	1.0	21
95-50-1	1,2-Dichlorobenzene	1.0	20
541-73-1	1,3-Dichlorobenzene	1.0	20
106-46-7	1,4-Dichlorobenzene	1.0	19
75-71-8	Dichlorodifluoromethane	1.0	20
75-34-3	1,1-Dichloroethane	1.0	20
107-06-2	1,2-Dichloroethane	1.0	20
75-35-4	1,1-Dichloroethene	1.0	19
156-59-2	cis-1,2-Dichloroethene	1.0	20
156-60-5	trans-1,2-Dichloroethene	1.0	19
78-87-5	1,2-Dichloropropane	1.0	20
10061-01-5	cis-1,3-Dichloropropene	1.0	20
10061-02-6	trans-1,3-Dichloropropene	1.0	20
75-09-2	Methylene chloride (Dichloromethane)	5.0	20
79-34-5	1,1,2,2-Tetrachlorethane	1.0	21
127-18-4	Tetrachloroethene	1.0	20
71-55-6	1,1,1-Trichloroethane	1.0	19
79-00-5	1,1,2-Trichloroethane	1.0	21
79-01-6	Trichloroethene	1.0	20
75-69-4	Trichlorofluoromethane	1.0	19
75-01-4	Vinyl chloride	1.0	- 20
	SURROGATE-Fluorobenzene (QC Limits - 61-133	3%)	100 % Rec

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333-PC4-1

## REPORT OF ANALYTICAL RESULTS PURGEABLE HALOCARBONS

Date collected: 7/11/96
Date extracted: N/A
Date analyzed: 7/19/96

Matrix: Water
Method: EPA601 (MOD)

% Moisture: 100

Sample Group: MB370

Lab Sample ID: MB370004RE Lab file 1 ID: G19T019 Lab file 2 ID: G19U019

Dilution factor: 1.0
Reporting units: ug/L

CAS NUMBER	COMPOUND NAME F	REPORTING LIMIT	RESULT
75-27-4	Bromodichloromethane	1.0	(
75-25-2	Bromoform	1.0	τ
74-83-9	Bromomethane	1.0	τ
56-23-5	Carbon tetrachloride	1.0	τ
108-90-7	Chlorobenzene	1.0	τ
75-00-3	Chloroethane	1.0	Ţ
110-75-8	2-Chloroethyl vinyl ether	1.0	τ
67-66-3	Chloroform	1.0	τ
74-87-3	Chloromethane	1.0	٠ ر
124-48-1	Dibromochloromethane	1.0	Ţ
95-50-1	1,2-Dichlorobenzene	1.0	Ţ
541-73-1	1,3-Dichlorobenzene	1.0	ı
106-46-7	1,4-Dichlorobenzene	1.0	Ţ
75-71-8	Dichlorodifluoromethane	1.0	ı
75-34-3	1,1-Dichloroethane	1.0	Ţ
107-06-2	1,2-Dichloroethane	1.0	Ţ
75-35-4	1,1-Dichloroethene	1.0	
156-59-2	cis-1,2-Dichloroethene	1.0	Ţ
156-60-5	trans-1,2-Dichloroethene	1.0	Ţ
78-87-5	1,2-Dichloropropane	1.0	Į
10061-01-5	cis-1,3-Dichloropropene	1.0	Ţ
10061-02-6	trans-1,3-Dichloropropene	1.0	ī
75-09-2	Methylene chloride (Dichloromethane)	5.0	Ţ
79-34-5	1,1,2,2-Tetrachlorethane	1.0	Ţ
127-18-4	Tetrachloroethene	1.0	1
71-55-6	1,1,1-Trichloroethane	1.0	1
79-00-5	1,1,2-Trichloroethane	1.0	1
79-01-6	Trichloroethene	1.0	1
75-69-4	Trichlorofluoromethane	1.0	1
75-01-4	Vinyl chloride	1.0	······
	SURROGATE-Fluorobenzene (QC Limits - 61-133	3%)	92 % Red

Jul

333-MW01-1B

## REPORT OF ANALYTICAL RESULTS PURGEABLE HALOCARBONS

 Date collected:
 7/11/96
 Sample Group:
 MB370

 Date extracted:
 N/A
 Lab Sample ID:
 MB370005

 Date analyzed:
 7/19/96
 Lab file 1 ID:
 G18T022

 Matrix:
 Water
 Lab file 2 ID:
 G18U022

 Method:
 EPA601 (MOD)
 Dilution factor:
 1.0

 % Moisture:
 100
 Reporting units:
 ug/L

CAS NUMBER	COMPOUND NAME	REPORTING LIMIT	RESULT
75-27-4	Bromodichloromethane	1.0	บ
75-25-2	Bromoform	1.0	บ
74-83-9	Bromomethane	1.0	U
56-23-5	Carbon tetrachloride	1.0	บ
108-90-7	Chlorobenzene	1.0	U
75-00-3	Chloroethane	1.0	U
110-75-8	2-Chloroethyl vinyl ether	1.0	U
67-66-3	Chloroform	1.0	บ
74-87-3	Chloromethane	1.0	U
124-48-1	Dibromochloromethane	1.0	U
95-50-1	1,2-Dichlorobenzene	1.0	U
541-73-1	1,3-Dichlorobenzene	1.0	t
106-46-7	1,4-Dichlorobenzene	1.0	τ
75-71-8	Dichlorodifluoromethane	1.0	τ
75-34-3	1,1-Dichloroethane	1.0	τ
107-06-2	1,2-Dichloroethane	1.0	U
75-35-4	1,1-Dichloroethene	1.0	Ţ
156-59-2	cis-1,2-Dichloroethene	1.0	į
156-60-5	trans-1,2-Dichloroethene	1.0	Ţ
78-87-5	1,2-Dichloropropane	1.0	τ
10061-01-5	cis-1,3-Dichloropropene	1.0	ι
10061-02-6	trans-1,3-Dichloropropene	1.0	τ
75-09-2	Methylene chloride (Dichloromethane) 5.0		J
79-34-5	1.1.2.2-Tetrachlorethane 1.0		Ţ
127-18-4	Tetrachloroethene	1.0	τ
71-55-6	1,1,1-Trichloroethane	1.0	Ţ
79-00-5	1,1,2-Trichloroethane	1.0	τ
79-01-6	Trichloroethene	1.0	Ţ
75-69-4	Trichlorofluoromethane	1.0	τ
75-01-4	Vinyl chloride	1.0	٦ ,
	SURROGATE-Fluorobenzene (QC Limits - 61-13	33%)	103 % Rec

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333-MW03-1D

## REPORT OF ANALYTICAL RESULTS PURGEABLE HALOCARBONS

 Date collected:
 7/11/96
 Sample Group:
 MB370

 Date extracted:
 N/A
 Lab Sample ID:
 MB370006

 Date analyzed:
 7/18/96
 Lab file 1 ID:
 G18T017

 Matrix:
 Water
 Lab file 2 ID:
 G18U017

 Method:
 EPA601 (MOD)
 Dilution factor:
 1.0

 % Moisture:
 100
 Reporting units:
 ug/L

CAS NUMBER	COMPOUND NAME F	REPORTING LIMIT	RESUL	T
75-27-4	Bromodichloromethane	1.0		บ
75-25-2	Bromoform	1.0		U
74-83-9	Bromomethane	1.0		U
56-23-5	Carbon tetrachloride	1.0		U
108-90-7	Chlorobenzene	1.0		U
75-00-3 .	Chloroethane	1.0		U
110-75-8	2-Chloroethyl vinyl ether	1.0		U
67-66-3	Chloroform	1.0		U
74-87-3	Chloromethane	1.0		U
124-48-1	Dibromochloromethane	1.0		U
95-50-1	1,2-Dichlorobenzene	1.0		τ
541-73-1	1,3-Dichlorobenzene	1.0		ι
106-46-7	1,4-Dichlorobenzene	1.0		τ
75-71-8	Dichlorodifluoromethane	1.0		ι
75-34-3	1,1-Dichloroethane	1.0		τ
107-06-2	1,2-Dichloroethane	1.0		τ
75-35-4	1,1-Dichloroethene	1.0		τ
156-59-2	cis-1,2-Dichloroethene	1.0		τ
156-60-5	trans-1,2-Dichloroethene	1.0		τ
78-87-5	1,2-Dichloropropane	1.0		τ
10061-01-5	cis-1,3-Dichloropropene	1.0		τ
10061-02-6	trans-1,3-Dichloropropene	1.0		ŧ
75-09-2	Methylene chloride (Dichloromethane)	• •		ξ
79-34-5	1,1,2,2-Tetrachlorethane	1.0		τ
127-18-4	Tetrachloroethene	1.0		Į
71-55-6	1,1,1-Trichloroethane	1.0		τ
79-00-5	1,1,2-Trichloroethane	1.0		ŧ
79-01-6	Trichloroethene	1.0		τ
75-69-4	Trichlorofluoromethane	1.0		τ
.75-01-4	Vinyl chloride	1.0	7	τ
	SURROGATE-Fluorobenzene (QC Limits - 61-133	3%)	102 % R	ec

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## REPORT OF ANALYTICAL RESULTS PURGEABLE HALOCARBONS

Date collected: 7/11/96
Date extracted: N/A
Date analyzed: 7/18/96

% Moisture: 100

Matrix: Water
Method: EPA601 (MOD)

Sample Group: MB370 Lab Sample ID: MB370007

Lab file 1 ID: G18T018
Lab file 2 ID: G18U018

Dilution factor: 1.0
Reporting units: ug/L

CAS NUMBER	COMPOUND NAME	REPORTING LIMIT	RESULT
75-27-4	Bromodichloromethane	1.0	
75-25-2	Bromoform	1.0	U
74-83-9	Bromomethane	1.0	U
56-23-5	Carbon tetrachloride	1.0	ប
108-90-7	Chlorobenzene	1.0	U
75-00-3	Chloroethane	1.0	U
110-75-8	2-Chloroethyl vinyl ether	1.0	U
67-66-3	Chloroform	1.0	U
74-87-3	Chloromethane	1.0	U
124-48-1	Dibromochloromethane	1.0	U
95-50-1	1,2-Dichlorobenzene	1.0	U
541-73-1 -	1,3-Dichlorobenzene	1.0	U
106-46-7	1,4-Dichlorobenzene	1.0	U
75-71-8	Dichlorodifluoromethane	1.0	U
75-34-3	1,1-Dichloroethane	1.0	U
107-06-2	1,2-Dichloroethane	1.0	U
75-35-4	1,1-Dichloroethene	1.0	U
156-59-2	cis-1,2-Dichloroethene	1.0	U
156-60-5	trans-1,2-Dichloroethene	1.0	U
78-87-5	1,2-Dichloropropane	1.0	U
10061-01-5	cis-1,3-Dichloropropene	1.0	U
10061-02-6	trans-1,3-Dichloropropene	1.0	U
75-09-2	Methylene chloride (Dichloromethane)	5.0	U
79-34-5	1,1,2,2-Tetrachlorethane	1.0	U
127-18-4	Tetrachloroethene	1.0	ប
71-55-6	1,1,1-Trichloroethane	1.0	υ
79-00-5	1,1,2-Trichloroethane	1.0	U
79-01-6	Trichloroethene	1.0	ប
75-69-4	Trichlorofluoromethane	1.0	U
75-01-4	Vinyl chloride	1.0	, U
	SURROGATE-Fluorobenzene (QC Limits - 61-13:	3%)	98 % Rec.

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# GC PURGEABLE AROMATICS

## CASE NARRATIVE GC PURGEABLE AROMATICS

QAL L	ab Refei	rence No./SDG. MB370
Proje	ct:	BROWN & ROOT COASTAL SYSTEMS STATION
ı.	RECEIPT	r
		eptions were encountered unless a Sample Receipt Exception Report is ed to the Chain-of-Custody included with this data package.
II.	HOLDING	G TIMES
	A. 8	Sample Preparation: Not applicable.
	В. 3	Sample Analysis: All holding times were met.
III:	METHOD	
	Cleanur	ation: N/A p: N/A is: EPA 602 (MOD)
IV.	PREPARI	ATION
	Not app	plicable.
v.	ANALYS	IS
	A. (	Calibration : All acceptance criteria were met.
	В. І	Blanks: All acceptance criteria were met.
	C. 8	Surrogates: All acceptance criteria were met.
	D.	Matrix Spikes: All acceptance criteria were met.
		Samples: Sample MB370001 (333-MW01-1) contained unidentified, non-target compounds.
	3	Primary analysis utilized a Restek Rtx 502.2 (105 meter x 0.53 mm) column. Confirmation analysis used a J&W . Scientific DB-VRX (75 meter x 0.45 mm) column.
agree excep hardc	d to by t for t opy data	at this data package is in compliance with the terms and conditions the client and QAL, Inc., both technically and for completeness he conditions noted above. Release of the data contained in this a package has been authorized by the Laboratory Manager or designated erified by the following signature.
SIGNE	for a	DATE: 08-06-9C  Merb Kelly Organic Division Manager

GC PURGEABLE AROMATICS
Lab Reference No./SDG: MB370
Page 2

# CASE NARRATIVE Addendum

Sample Information

LAB SAMPLE ID	CLIENT SAMPLE ID	SAMPLE MATRIX	DATE SAMPLED	DATE EXTRACTED	DATE ANALYZED	SAMPLE pH <sup>1</sup>
MB370001	333-MW01-1	WATER	7/11/96	N/A	7/19/96	<2
MB370002RE	333-MW02-1	WATER	7/11/96	N/A	7/19/96	<2
MB370003	333-MW03-1	WATER	7/11/96	N/A	7/18/96	<2
MB370003MS	333-MW03-1MS	WATER	7/11/96	N/A	7/19/96	<2
MB370003MSD	333-MW03-1MD	WATER	7/11/96	N/A	7/19/96	<2
MB370004RE	333-PC4-1	WATER	7/11/96	N/A	7/19/96	<2
MB370005	333-MW01-1B	WATER	7/11/96	N/A	7/19/96	<2
MB370006	333-MW03-1D	WATER	7/11/96	N/A	7/18/96	<2
MB370007	TRIP BLANK	WATER	7/11/96	N/A	7/18/96	<2
WMV096G182	VBLK001	WATER	N/A	N/A	7/18/96	N/A
WMV096G191	VBLK002	WATER	N/A	N/A	7/19/96	N/A

<sup>1</sup> Applies to samples designated for purgeable VOA analysis only.

#### CURRENT METHOD DETECTION LIMITS (MDLs) PURGEABLE AROMATICS

Date collected: N/A

Sample Group: LABQC Lab Sample ID: Multiple Samples Date extracted: N/A

Date analyzed: 3/13/96 Lab file 1 ID: N/A Matrix: Water Lab file 2 ID: N/A Method: EPA602 (MOD) Dilution factor: 1.0 % Moisture: 100 Reporting units: ug/L

CAS NUMBER	COMPOUND NAME	REPORTING LIMIT	RESULT
71-43-2	Benzene	1.0	0.102
108-90-7	Chlorobenzene	1.0	0.141
95-50-1	1.2-Dichlorobenzene	1.0	0.128
541-73-1	1,3-Dichlorobenzene	1.0	0.137
106-46-7	1,4-Dichlorobenzene	1.0	0.145
100-41-4	Ethylbenzene	1.0	0.129
1634-04-4	tert-butyl methyl ether	1.0	0.087
108-88-3	Toluene	1.0	0.102
108-38-3/106-42-3	m-, p-Xylene	2.0	0.312
95-47-6	o-Xylene	1.0	0.189

333-MW01-1

## REPORT OF ANALYTICAL RESULTS PURGEABLE AROMATICS

 Date collected:
 7/11/96
 Sample Group:
 MB370

 Date extracted:
 N/A
 Lab Sample ID:
 MB370001

 Date analyzed:
 7/19/96
 Lab file 1 ID:
 G18T019

 Matrix:
 Water
 Lab file 2 ID:
 G18U019

 Method:
 EPA602 (MOD)
 Dilution factor:
 1.0

 % Moisture:
 100
 Reporting units:
 ug/L

CAS NUMBER	COMPOUND NAME	REPORTING LIMIT	RESULT
71-43-2	Benzene	1.0	U
108-88-3	Toluene	1.0	1.3
100-41-4	Ethylbenzene	1.0	6.7
1330-20-7	Xylenes (Total)	1.0	6.8
N/A	Total Volatile Organic Aromatics	1.0	15
1634-04-4	Methyl-tert-butyl ether	1.0	9.0
	SURROGATE-Fluorobenzene (QC Limits - 61-1)	33%)	101 % Rec.

Jul

333-MW02-1

#### REPORT OF ANALYTICAL RESULTS PURGEABLE AROMATICS

Date collected: 7/11/96 Sample Group: MB370 Lab Sample ID: MB370002RE Date extracted: N/A Date analyzed: 7/19/96 Lab file 1 ID: G19T017 Matrix: Water Lab file 2 ID: G19U017 Method: EPA602 (MOD) Dilution factor: 1.0

% Moisture: 100

Reporting units: ug/L

CAS. NUMBER	COMPOUND NAME	REPORTING LIMIT	RESULT
71-43-2	Benzene	1.0	τ
108-88-3	Toluene	1.0	τ
100-41-4	Ethylbenzene	1.0	ι
1330-20-7	Xylenes (Total)	1.0	τ
N/A	Total Volatile Organic Aromatics	1.0	Ţ
1634-04-4	Methyl-tert-butyl ether	1.0	τ
	SURROGATE-Fluorobenzene (QC Limits -	· 61-133%)	99 % Rec

333-MW03-1

## REPORT OF ANALYTICAL RESULTS PURGEABLE AROMATICS

Date collected: 7/11/96
Date extracted: N/A
Date analyzed: 7/18/96
Matrix: Water
Method: EPA602 (MOD)

% Moisture: 100

Sample Group: MB370
Lab Sample ID: MB370003
Lab file 1 ID: G18T016
Lab file 2 ID: G18U016
Dilution factor: 1.0
Reporting units: ug/L

CAS NUMBER	COMPOUND NAME	REPORTING LIMIT	RESULT
71-43-2	Benzene	1.0	บ
108-88-3	Toluene	1.0	U
100-41-4	Ethylbenzene	1.0	U
1330-20-7	Xylenes (Total)	1.0	U
N/A	Total Volatile Organic Aromatics	1.0	U
1634-04-4	Methyl-tert-butyl ether	1.0	ប
	SURROGATE-Fluorobenzene (QC Limits - 6	1-133%)	99 % Rec.

Jul

333-MW03-1MS

## REPORT OF ANALYTICAL RESULTS PURGEABLE AROMATICS

CAS NUMBER	COMPOUND NAME	REPORTING LIMIT	RESULT
71-43-2	Benzene	1.0	20
108-88-3	Toluene	1.0	20
100-41-4	Ethylbenzene	1.0	20
1330-20-7	Xylenes (Total)	1.0	60
N/A	Total Volatile Organic Aromatics	1.0	N/A
1634-04-4	Methyl-tert-butyl ether	1.0	23
	SURROGATE-Fluorobenzene (QC Limits - 61-	-133%)	100 % Rec.



333-MW03-1MD

## REPORT OF ANALYTICAL RESULTS PURGEABLE AROMATICS

 Date collected:
 7/11/96
 Sample Group:
 MB370

 Date extracted:
 N/A
 Lab Sample ID:
 MB370003MSD

 Date analyzed:
 7/19/96
 Lab file 1 ID:
 G18T027

 Matrix:
 Water
 Lab file 2 ID:
 G18U027

 Method:
 EPA602 (MOD)
 Dilution factor:
 1.0

 % Moisture:
 100
 Reporting units:
 ug/L

CAS NUMBER COMPOUND NAME REPORTING LIMIT RESULT

71-43-2 Benzene 1.0 21
108-88-3 Toluene 1.0 21

71-43-2 108-88-3 Toluene 1.0 21 100-41-4 Ethylbenzene 1.0 20 1330-20-7 Xylenes (Total) 1.0 61 N/A Total Volatile Organic Aromatics 1.0 N/A 1634-04-4 Methyl-tert-butyl ether 1.0 24 SURROGATE-Fluorobenzene (QC Limits - 61-133%) 100 % Rec.

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#### CLIENT SAMPLE ID

333-PC4-1

#### REPORT OF ANALYTICAL RESULTS PURGEABLE AROMATICS

Date collected: 7/11/96 Date extracted: N/A Date analyzed: 7/19/96

Matrix: Water Method: EPA602 (MOD)

% Moisture: 100

Sample Group: MB370 Lab Sample ID: MB370004RE Lab file 1 ID: G19T019

Lab file 2 ID: G19U019 Dilution factor: 1.0

Reporting units: ug/L

CAS NUMBER	COMPOUND NAME	REPORTING LIMIT	RESULT
71-43-2	Benzene	1.0	U
108-88-3	Toluene	1.0	U
100-41-4	Ethylbenzene	1.0	U
1330-20-7	Xylenes (Total)	1.0	U
N/A	Total Volatile Organic Aromatics	1.0	Ū
1634-04-4	Methyl-tert-butyl ether	1.0	ប
	SURROGATE-Fluorobenzene (QC Limits - 61-13	3%)	92 % Rec.

333-MW01-1B

## REPORT OF ANALYTICAL RESULTS PURGEABLE AROMATICS

 Date collected:
 7/11/96
 Sample Group:
 MB370

 Date extracted:
 N/A
 Lab Sample ID:
 MB370005

 Date analyzed:
 7/19/96
 Lab file 1 ID:
 G18T022

 Matrix:
 Water
 Lab file 2 ID:
 G18U022

 Method:
 EPA602 (MOD)
 Dilution factor:
 1.0

 % Moisture:
 100
 Reporting units:
 ug/L

CAS NUMBER	COMPOUND NAME	REPORTING LIMIT	RESULT
71-43-2	Benzene	1.0	U
108-88-3	Toluene	1.0	1.4
100-41-4	Ethylbenzene	1.0	ប
1330-20-7	Xylenes (Total)	1.0	U
N/A	Total Volatile Organic Aromatics	1.0	1.4
1634-04-4	Methyl-tert-butyl ether	1.0	ប
	SURROGATE-Fluorobenzene (QC Limits - 61-13	3%)	103 % Rec.



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## REPORT OF ANALYTICAL RESULTS PURGEABLE AROMATICS

 Date collected:
 7/11/96
 Sample Group:
 MB370

 Date extracted:
 N/A
 Lab Sample ID:
 MB370007

 Date analyzed:
 7/18/96
 Lab file 1 ID:
 G18T018

 Matrix:
 Water
 Lab file 2 ID:
 G18U018

 Method:
 EPA602 (MOD)
 Dilution factor:
 1.0

 % Moisture:
 100
 Reporting units:
 ug/L

CAS NUMBER	COMPOUND NAME	REPORTING LIMIT	RESULT
71-43-2	Benzene	1.0	ָּ
108-88-3	Toluene	1.0	1.8
100-41-4	Ethylbenzene	1.0	U
1330-20-7	Xylenes (Total)	1.0	บ
N/A	Total Volatile Organic Aromatics	1.0	1.8
1634-04-4	Methyl-tert-butyl ether	1.0	U
	SURROGATE-Fluorobenzene (QC Limits - 61-	133%)	98 % Rec.



# GC EXTRACTABLE VOLATILE ORGANICS (EDB)

## CASE NARRATIVE GC EXTRACTABLE VOLATILE ORGANICS (EDB)

QAL Lab	Reference	No./SDC	. <u>MB370</u>	_
Project	Brown &	Root Co	astal Systems Station	

#### I. RECEIPT

No exceptions were encountered unless a Sample Receipt Exception Report is attached to the Chain-of-Custody included with this data package.

#### II. HOLDING TIMES

- A. Sample Preparation: All holding times were met.
- B. Sample Analysis: All holding times were met.

#### III. METHOD

Preparation: N/A Cleanup: N/A Analysis: EPA 504.1

#### IV. PREPARATION

Sample preparation proceeded normally.

#### V. ANALYSIS

- A. Calibration: All acceptance criteria were met.
- B. Blanks: All acceptance criteria were met.
- C. Surrogates: All acceptance criteria were met.
- D. Spikes: Water matrix spikes were performed using a sample from this contract. The summary of the MS, MSD, and the associated blank spike results have been included in this data package.
- E. Samples: Sample analysis proceeded normally.

A summary of current applicable method detection limits (MDLs) immediately follows this case narrative.

GC EXTRACTABLE VOLATILE ORGANICS (EDB) Lab Reference No./SDG: MB370

Page 2

I certify that this data package is in compliance with the terms and conditions agreed to by the client and QAL, Inc., both technically and for completeness except for the conditions noted above. Release of the data contained in this hardcopy data package has been authorized by the Laboratory Manager or designated person, as verified by the following signature.

SIGNED: ~

\_ DATE: 7/31/96

GC EXTRACTABLE VOLATILE ORGANICS (EDB)
Lab Reference No./SDG: MB370
Page 3

# CASE NARRATIVE Addendum

### Sample Information

LAB	CLIENT	SAMPLE	DATE	DATE	DATE	SAMPLE
SAMPLE ID	SAMPLE ID	MATRIX	SAMPLED	EXTRACTED	ANALYZED	DH <sub>T</sub>
MB370001	333-MW01-1	WATER	07/11/96	07/29/96	07/29/96	N/A
MB370002	333-MW02-1	WATER	07/11/96	07/29/96	07/29/96	N/A
MB370003	333-MW03-1	WATER	07/11/96	07/29/96	07/29/96	N/A
MB370003MS	333-MW03-1MS	WATER	07/11/96	07/29/96	07/29/96	N/A
MB370003MSD	333-MW03-1MD	WATER	07/11/96	07/29/96	07/29/96	N/A
MB370004	333-PC4-1	WATER	07/11/96	07/29/96	07/29/96	N/A
MB370005	333-MW01-1B	WATER	07/11/96	07/29/96	07/29/96	N/A
MB370006	333-MW03-1D	WATER	07/11/96	07/29/96	07/29/96	N/A
W07296B1	OC BLANK	WATER	N/A	07/29/96	07/29/96	N/A

<sup>1</sup> Applies to samples designated for purgeable VOA analysis only.

#### ORGANICS ANALYSIS METHOD DETECTION LIMITS

#### GC EXTRACTABLE VOLATILE ORGANICS (EDB)

Laboratory Name: <u>CH2M HILL</u> Sample Matrix: <u>WATER</u>

Analytical Method: 504.1

CAS Number Compound ug/L

106-93-4 1,2-Dibromoethane (EDB) 0.003

Aboratory Name: CH2M HILL Concentration: LOW Date Extracted: 07/29/96

Lab Sample ID: MB370001 Sample Matrix: WATER Date Analyzed: 07/29/96

Client Sample ID: 333-MW01-1 Percent Moisture: Dilution Factor: 1.0

#### GC EXTRACTABLE VOLATILE ORGANICS (EDB)

CAS Number uq/L
106-93-4 1,2-Dibromoethane (EDB) . . . 0.02 U

1,1,2,2-Tetrachloroethane - SS 96

U - Analyzed for but not detected.

B - Detected in QC blank.

J - Detected, concentration estimated.

SS - Surrogate Standard reported as percent recovery.

Comments:

Form I

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aboratory Name: CH2M HILL Concentration: LOW Date Extracted: 07/29/96
Lab Sample ID: MB370002 Sample Matrix: WATER Date Analyzed: 07/29/96
Client Sample ID: 333-MW02-1 Percent Moisture: Dilution Factor: 1.0

#### GC EXTRACTABLE VOLATILE ORGANICS (EDB)

- U Analyzed for but not detected.
- B Detected in QC blank.
- J Detected, concentration estimated.
- SS Surrogate Standard reported as percent recovery.

Comments:

Form I

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Aboratory Name: CH2M HILL Concentration: LOW Date Extracted: 07/29/96

Lab Sample ID: MB370003 Sample Matrix: WATER Date Analyzed: 07/29/96

Client Sample ID: 333-MW03-1 Percent Moisture: Dilution Factor: 1.0

#### GC EXTRACTABLE VOLATILE ORGANICS (EDB)

U - Analyzed for but not detected.

B - Detected in QC blank.

J - Detected, concentration estimated.

SS - Surrogate Standard reported as percent recovery.

Comments:

Form I

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aboratory Name: CH2M HILL Concentration: LOW Date Extracted: 07/29/96
Lab Sample ID: MB370004 Sample Matrix: WATER Date Analyzed: 07/29/96
Client Sample ID: 333-PC4-1 Percent Moisture: Dilution Factor: 1.0

#### GC EXTRACTABLE VOLATILE ORGANICS (EDB)

CAS Number uq/L
106-93-4 1,2-Dibromoethane (EDB) . . . 0.02 U

1,1,2,2-Tetrachloroethane - SS 98

U - Analyzed for but not detected.

B - Detected in QC blank.

J - Detected, concentration estimated.

SS - Surrogate Standard reported as percent recovery.

Comments:

Form I

A NO

aboratory Name: CH2M HILL Concentration: LOW ·Date Extracted: 07/29/96 Lab Sample ID: MB370005 Sample Matrix: Date Analyzed: WATER 07/29/96 Client Sample ID: 333-MW01-1B Percent Moisture: Dilution Factor: \_\_\_\_1.0

#### GC EXTRACTABLE VOLATILE ORGANICS (EDB)

CAS Number uq/L 106-93-4 1,2-Dibromoethane (EDB) . . . 0.02 U 1,1,2,2-Tetrachloroethane - SS 95

U - Analyzed for but not detected.

B - Detected in QC blank.

J - Detected, concentration estimated.

SS - Surrogate Standard reported as percent recovery.

Comments:

Form I

aboratory Name: CH2M HILL Concentration: LOW .Date Extracted: 07/29/96

Lab Sample ID: MB370006 Sample Matrix: WATER Date Analyzed: 07/29/96

Client Sample ID: 333-MW03-1D Percent Moisture: Dilution Factor: 1.0

#### GC EXTRACTABLE VOLATILE ORGANICS (EDB)

CAS Number ug/L

106-93-4 1,2-Dibromoethane (EDB) . . . 0.02 U

1,1,2,2-Tetrachloroethane - SS 95

- U Analyzed for but not detected.
- B Detected in QC blank.
- J Detected, concentration estimated.
- SS Surrogate Standard reported as percent recovery.

Comments:

Form I

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(334) 271-2440

GC POLYNUCLEAR AROMATIC HYDROCARBONS

## CASE NARRATIVE GC POLYNUCLEAR AROMATIC HYDROCARBONS

QAL	Lab	Reference	No./SDG.	<u>MB370</u>
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Project: Brown & Root Coastal Systems Station

#### I. RECEIPT

No exceptions were encountered unless a Sample Receipt Exception Report is attached to the Chain-of-Custody included with this data package.

#### II. HOLDING TIMES

- A. Sample Preparation: All holding times were met.
- B. Sample Analysis: All holding times were met.

#### III. METHOD

Preparation: N/A Cleanup: N/A Analysis: EPA 610

#### IV. PREPARATION

Sample preparation proceeded normally.

#### V. ANALYSIS

A. Calibration: All acceptance criteria were met.

Both the initial calibration and continuing calibration summaries include data for both the primary and confirmation columns. Each compound will appear in the summary reports twice. The first time the compound will not be proceeded by the "#" symbol, referring to compounds identified from the first column (RTX-5); the next time it will have the "#"symbol, referring to compounds identified from the second column (RTX-200) (for example, Naphthalene and #Naphthalene).

- B. Blanks: All acceptance criteria were met.
- C. Surrogates: All acceptance criteria were met.
- D. Spikes: Matrix spikes were performed using a sample from this contract. The summary of the MS/MSD results has been included in this data package.
- E. Samples: Sample analysis proceeded normally.

A summary of current applicable method detection limits (MDLs) immediately follows this case narrative.

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GC	POLYNUCLEAR	R AROMATIC	HYDROCARBONS	
Lab	Reference	No./SDG:	MB370	
Pag	re 2			

F. Other: Primary and confirmation data were simultaneously acquired using two dissimilar analytical columns (RTX-5 and RTX-200) connected in parallel to one injection port and one detector.

DATE: 7/31/96

I certify that this data package is in compliance with the terms and conditions agreed to by the client and QAL, Inc., both technically and for completeness except for the conditions noted above. Release of the data contained in this hardcopy data package has been authorized by the Laboratory Manager or designated person, as verified by the following signature.

SIGNED / Anum
Tammy Carey

Chemist III

GC POLYNUCLEAR AROMATIC HYDROCARBONS
Lab Reference No./SDG: MB370
Page 3

# CASE NARRATIVE Addendum

Sample Information

LAB	CLIENT	SAMPLE	DATE	DATE	DATE ANALYZED	SAMPLE
SAMPLE ID	SAMPLE ID	MATRIX	SAMPLED	EXTRACTED		pH1
MB370001	333-MW01-1	WATER	07/11/96	07/15/96	07/25/96	N/A
MB370002	333-MW02-1	WATER	07/11/96	07/15/96	07/25/96	N/A
MB370003	333-MW03-1	WATER	07/11/96	07/15/96	07/25/96	N/A
MB370003MS	333-MW03-1MS	WATER	07/11/96	07/17/96	07/25/96	N/A
MB370003MSD	333-MW03-1MD	WATER	07/11/96	07/17/96	07/25/96	N/A
MB370004	333-PC4-1	WATER	07/11/96	07/15/96	07/25/96	N/A
MB370005	333-MW01-1B	WATER	07/11/96	07/15/96	07/25/96	N/A
MB370006	333-MW03-1D	WATER	07/11/96	07/15/96	07/25/96	N/A
W07156B1	NBLK06	WATER	N/A	07/15/96	07/24/96	N/A
W07176B1	NBLK07	WATER	n/a	07/17/96	07/25/96	N/A

<sup>1</sup> Applies to samples designated for purgeable VOA analysis only.

### ORGANICS ANALYSIS METHOD DETECTION LIMITS

### POLYNUCLEAR AROMATIC HYDROCARBON (PNA) COMPOUNDS

Laboratory Name: CH2M HILL Sample Matrix: WATER

Analytical Method: 610 Extraction: SEPARATORY FUNNEL

Compound	MDL ug/L
Compound	<u>uq/ บ</u>
Naphthalene	0.57
2-Methylnaphthalene	0.47
1-Methylnaphthalene	0.42
Acenaphthylene	0.38
Acenaphthene	0.41
Fluorene	0.38
Phenanthrene	0.51
Anthracene	0.32
Fluoranthene	0.60
Pyrene	0.27
Benzo(a)anthracene	0.24
Chrysene	0.49
Benzo(b)fluoranthene	0.40
Benzo(k)fluoranthene	0.39
Benzo(a)pyrene	0.33
Indeno (123-cd) pyrene	0.40
Dibenzo(ah)anthracene	0.30
Benzo(ghi)perylene	0.27

Laboratory Name: CH2M HILL Concentration: LOW Date Extracted: 07/15/96
Lab Sample ID: MB370001 Sample Matrix: WATER Date Analyzed: 07/25/96
Client Sample ID: 333-MW01-1 Volume Extracted: 1005mL Dilution Factor: 1.0

#### PNA COMPOUNDS

CAS Number		ug/L	
91-20-3	Naphthalene	2	U
91-57-6	2-Methylnaphthalene	2	U
90-12-0	1-Methylnaphthalene	2	U
208-96-8	Acenaphthylene	2	U
83-32-9	Acenaphthene	2	Ū
86-73-7	Fluorene	2	U
85-01-8	Phenanthrene	2	U
120-12-7	Anthracene	2	U
206-44-0	Fluoranthene	2	U
129-00-0	Pyrene	2	υ
56-55-3	Benzo(a) anthracene	2	Ŭ
218-01-9	Chrysene	2	U
205-99-2	Benzo(b) fluoranthene	2	U
207-08-9	Benzo(k) fluoranthene	2	σ
50-32-8	Benzo(a)pyrene	2	U
193-39-5	<pre>Indeno(1,2,3-cd)pyrene</pre>	2	Ū
53-70-3	Dibenzo(a,h)anthracene	2	U
191-24-2	Benzo(g,h,i)perylene	2	σ
	Terphenyl-d14 - SS	58	 8

- U Analyzed for but not detected.
- B Detected in QC blank.
- J Detected, concentration estimated.
- SS Surrogate Standard reported as percent recovery.

## Comments:

Form I

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Laboratory Name: CH2M HILL Concentration: LOW Date Extracted: 07/15/96
Lab Sample ID: MB370002 Sample Matrix: WATER Date Analyzed: 07/25/96
Client Sample ID: 333-MW02-1 Volume Extracted: 1055mL Dilution Factor: 1.0

### PNA COMPOUNDS

CAS Number		ug/L	
91-20-3	Naphthalene	2	U
91-57-6	2-Methylnaphthalene	2	U
90-12-0	1-Methylnaphthalene	2	σ
208-96-8	Acenaphthylene	2	U
83-32-9	Acenaphthene	2	U
86-73-7	Fluorene	2	U
85-01-8	Phenanthrene	2	U
120-12-7	Anthracene	2	U
206-44-0	Fluoranthene	2	U
129-00-0	Pyrene	2	U
56-55-3	Benzo(a) anthracene	2	Ŭ
218-01-9	Chrysene	2	Ū
205-99-2	Benzo(b) fluoranthene	2	U
207-08-9	Benzo(k)fluoranthene	2	ט
50-32-8	Benzo(a)pyrene	2	U
193-39-5	Indeno(1,2,3-cd)pyrene	2	U
53-70-3	Dibenzo(a,h)anthracene	2	Ū
191-24-2	Benzo(g,h,i)perylene	2	ָּט
	Terphenyl-d14 - SS	72	%

#### Comments:

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U - Analyzed for but not detected.

B - Detected in QC blank.

J - Detected, concentration estimated.

SS - Surrogate Standard reported as percent recovery.

Laboratory Name: CH2M HILL Concentration: LOW Date Extracted: 07/15/96
Lab Sample ID: MB370003 Sample Matrix: WATER Date Analyzed: 07/25/96
Client Sample ID: 333-MW03-1 Volume Extracted: 1065mL Dilution Factor: 1.0

#### PNA COMPOUNDS

CAS Number		uq/L	
91-20-3	Naphthalene	2	ש
91-57-6	2-Methylnaphthalene	2	U
90-12-0	1-Methylnaphthalene	2	U
208-96-8	Acenaphthylene	2	U
83-32-9	Acenaphthene	2	Ū
86-73-7	Fluorene	2	Ū
85-01-8	Phenanthrene	2	U
120-12-7	Anthracene	2	U
206-44-0	Fluoranthene	2	U
129-00-0	Pyrene	2	σ
56-55-3	Benzo(a) anthracene	2	U
218-01-9	Chrysene	2	U
205-99-2	Benzo(b) fluoranthene	2	U
207-08-9	Benzo(k) fluoranthene	2	σ
50-32-8	Benzo(a)pyrene	2	U
193-39-5	Indeno(1,2,3-cd)pyrene	2	Ū
53-70-3	Dibenzo(a,h)anthracene	2	Ū
191-24-2	Benzo(g,h,i)perylene	2	U
	Terphenyl-d14 - SS	72	ક

## Comments:

Form I

U - Analyzed for but not detected.

B-- Detected in QC blank.

J - Detected, concentration estimated.

SS - Surrogate Standard reported as percent recovery.

Laboratory Name: CH2M HILL Concentration: LOW Date Extracted: 07/15/96
Lab Sample ID: MB370004 Sample Matrix: WATER Date Analyzed: 07/25/96
Client Sample ID: 333-PC4-1 Volume Extracted: 1020mL Dilution Factor: \_\_\_\_\_1.0

#### PNA COMPOUNDS

CAS Number		ug/L	
91-20-3	Naphthalene	2	Ū
91-57-6	2-Methylnaphthalene	2	U
90-12-0	1-Methylnaphthalene	2	U
208-96-8	Acenaphthylene	2	U
83-32-9	Acenaphthene	2	U
86-73-7	Fluorene	2	υ
85-01-8	Phenanthrene	2	U
120-12-7	Anthracene	2	U
206-44-0	Fluoranthene	2	ΰ
129-00-0	Pyrene	2	ΰ
56-55-3	Benzo(a) anthracene	2	U
218-01-9	Chrysene	2	U
205-99-2	Benzo(b) fluoranthene	2	U
207-08-9	Benzo(k) fluoranthene	2	ΰ
50-32-8	Benzo(a)pyrene	2	U
193-39-5	Indeno(1,2,3-cd)pyrene	2	U
53-70-3	Dibenzo(a,h)anthracene	2	U
191-24-2	Benzo(g,h,i)perylene	2	Ū
	Terphenyl-d14 - SS	74	ક

## Comments:

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U - Analyzed for but not detected.

B - Detected in QC blank.

J - Detected, concentration estimated.

SS - Surrogate Standard reported as percent recovery.

Laboratory Name: CH2M HILL Concentration: LOW Date Extracted: 07/15/96
Lab Sample ID: MB370005 Sample Matrix: WATER Date Analyzed: 07/25/96
Client Sample ID: 333-MW01-1B Volume Extracted: 1045mL Dilution Factor: 1.0

#### PNA COMPOUNDS

CAS Number		uq/L	
91-20-3	Naphthalene	2	Ū
91-57-6	2-Methylnaphthalene	2	U
90-12-0	1-Methylnaphthalene	2	U
208-96-8	Acenaphthylene	2	U
83-32-9	Acenaphthene	2	Ū
86-73-7	Fluorene	2	Ū
85-01-8	Phenanthrene	2	U
120-12-7	Anthracene	2	Ū
206-44-0	Fluoranthene	2	Ū
129-00-0	Pyrene	2	U
56-55-3	Benzo(a)anthracene	2	U
218-01-9	Chrysene	2	Ū
205-99-2	Benzo(b) fluoranthene	2	U
207-08-9	Benzo(k)fluoranthene	2	Ū
50-32-8	Benzo(a)pyrene	2	Ū
193-39-5	Indeno(1,2,3-cd)pyrene	2	Ū
53-70-3	Dibenzo(a,h)anthracene	2	σ
191-24-2	Benzo(g,h,i)perylene	2	<u>.</u>
	Terphenyl-d14 - SS	70	ક

### Comments:

Form I

U - Analyzed for but not detected.

B ~ Detected in QC blank.

J - Detected, concentration estimated.

SS - Surrogate Standard reported as percent recovery.

Laboratory Name: CH2M HILL Concentration: LOW Date Extracted: 07/15/96
Lab Sample ID: MB370006 Sample Matrix: WATER Date Analyzed: 07/25/96
Tlient Sample ID: 333-MW03-1D Volume Extracted: 1025mL Dilution Factor: 1.0

#### PNA COMPOUNDS

CAS Number		ug/L	
91-20-3	Naphthalene	2	U
91-57-6	2-Methylnaphthalene	2	U
90-12-0	1-Methylnaphthalene	2	U
208-96-8	Acenaphthylene	2	σ
83-32-9	Acenaphthene	2	Ū
86-73-7	Fluorene	2	U
85-01-8	Phenanthrene	2	U
120-12-7	Anthracene	2	σ
206-44-0	Fluoranthene	2	σ
129-00-0	Pyrene	2	U
56-55-3	Benzo(a) anthracene	2	Ū
218-01-9	Chrysene	2	U
205-99-2	Benzo(b) fluoranthene	2	Ū
207-08-9	Benzo(k) fluoranthene	2	υ
50-32-8	Benzo(a)pyrene	2	υ
193-39-5	Indeno(1,2,3-cd)pyrene	2	σ
53-70-3	Dibenzo(a,h)anthracene	2	σ
191-24-2	Benzo(g,h,i)perylene	2	Ū
	Terphenyl-d14 - SS	75	옿

## Comments:

U - Analyzed for but not detected.

B -- Detected in QC blank.

J - Detected, concentration estimated.

SS - Surrogate Standard reported as percent recovery.

CATIONS DATA PACKAGE

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## CASE NARRATIVE Cations

Labora	atory:	CH2M HILL Lab Ref. No.: MB370
Client	t/Proje	ect: Brown & Root Coastal Systems Station
I.		ng Time: olding times were met.
II.	Digest	tion Exceptions:
III.	Analys	sis:
	Α.	<pre>Calibration: All acceptance criteria were met.</pre>
	B.	Blanks: All acceptance criteria were met.
	C.	ICP Interference Check Sample: All acceptance criteria were met.
	D.	<pre>Spike Sample(s): All acceptance criteria were met.</pre>
	E.	<u>Duplicate Sample(s)</u> : All acceptance criteria were met.
	F.	<u>Laboratory Control Sample(s)</u> : All acceptance criteria were met.
	G.	ICP Serial Dilution: All acceptance criteria were met.
	H.	Other: None.
IV.	Any r	pt Exceptions:  eceipt exception will be addressed in a Sample Receipt Exception t which will be attached to the Chain-of-Custody in this package.
v.	Documo None.	entation Exceptions:
VI.	condi Inc., detai packa	tify that this data package is in compliance with the terms and tions agreed to by the client and Quality Analytical Laboratories, both technically and for completeness, except for the conditions led above. Release of the data contained in this hardcopy data ge has been authorized by the Laboratory Manager or his designee, rified by the following signature.  DATE: 1/3/96  Kaye Walker Inorganic Division Manager

## U.S. EPA - CLP

## 1 INORGANIC ANALYSES DATA SHEET

EPA SAMPLE NO.

		111011011111	ANALYSES DATA S	71111	_	
ab Name: CH2M	HILL		Contract: MI	B370	•	333-MW01-1
				•		SDG No.: MB370_
Matrix (soil/w	ater): WATE	R		Lab	Samp	le ID: MB370001
Level (low/med	): LOW_	_		Dat	e Rec	eived: 07/13/96
Solids:		0				
Co	ncentration	Units (ug	/L or mg/kg dry	y we	ight)	: UG/L_
	CAS No.	Analyte	Concentration	С	Q	м
	7440-43-9 7440-47-3 7439-92-1	Chromium_ Lead	2.9 3.3 21.4 16.9			F_P_P_F_
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Comments:						

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			ANALYSES DATA S		
ab Name: CH2M	HILL		Contract: Mi	3370	333-MW02-1
				MB370_	SDG No.: MB370
atrix (soil/w	ater): WATE	R		Lab Samp	le ID: MB370002_
evel (low/med	): LOW_	·-		Date Rec	eived: 07/13/96
Solids:		0			
Со	ncentration	Units (ug	/L or mg/kg dry	y weight)	: UG/L_
	CAS No.	Analyte	Concentration	c Q	M
	7440-38-2 7440-43-9 7440-47-3 7439-92-1	Arsenic Cadmium_ Chromium_ Lead	3.7 3.3 11.6 142	ש	F_P_P_F_
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olor After:	YELLOW	Clari	ty After: CLE	AR_	Artifacts:
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EPA SAMPLE NO.

Lab Name: CH2M	_HILL		Contract: M	337		333-MW03-1
						SDG No.: MB370_
Matrix (soil/w	ater): WATE	R		La	b Samp	ole ID: MB370003
Level (low/med	.): LOW_	_		Da	ite Rec	eived: 07/13/96
% Solids:		0				
Co	ncentration	Units (ug	/L or mg/kg dry	Z W	reight)	: UG/L_
	CAS No.	Analyte	Concentration	С	Q	M
	7440-38-2	Arsenic	2.6	핆		F_
		Cadmium	3.3			P_
	7440-47-3					[ P ]
	7439-92-1		5.8	_		[  F_
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Lab Code: MGM Case No.: MB370_ SAS No.: MB370_ SDG No.: MB370_  Matrix (soil/water): WATER	Lab Name: CH2M	HILL		Contract: MF	3370	333-PC4-1
Matrix (soil/water): WATER	•					
Level (low/med): LOW	Lab Code: MGM_	Cas	se No.: MB	370_ SAS No.:	: MB370_	SDG No.: MB370_
Concentration Units (ug/L or mg/kg dry weight): UG/L    CAS No.	Matrix (soil/wa	ater): WATE	R		Lab Samp	le ID: MB370004
Concentration Units (ug/L or mg/kg dry weight): UG/L_    CAS No.	Level (low/med	): LOW_	_		Date Rec	eived: 07/13/96
CAS No. Analyte Concentration C Q M    7440-38-2   Arsenic	% Solids:		0			
T440-38-2	Co	ncentration	Units (ug	/L or mg/kg dry	y weight)	: UG/L_
7440-43-9		CAS No.	Analyte	Concentration	C Q	м
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FORM I - IN

EPA SAMPLE NO.

Lab Name: CH2M	HILL	<del> </del>	Contract: M	B37		333-MW01-1B
Lab Code: MGM_	Ca	se No.: MB	370_ SAS No.:	: N	1B370_	SDG No.: MB370_
Matrix (soil/w	ater): WATE	R		La	ab Samp	le ID: MB370005
Level (low/med	): LOW_	_		Da	ate Rec	eived: 07/13/96
% Solids:		0				
Co	ncentration	Units (ug	/L or mg/kg dry	y v	veight)	: UG/L_
	CAS No.	Analyte	Concentration	С	Q	М
	7440-38-2	Arsenic	0.95	∣ਜ਼		F
		Cadmium	3.3		<del></del>	F_ P_ P_ F_
		Chromium	2.2			P
		Lead	0.75			F
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## U.S. EPA - CLP

## 1 INORGANIC ANALYSES DATA SHEET

EPA SAMPLE NO.

Lab Name: CH2M	HILL		Contract: Mi	B370	333-MW03-1D
					SDG No.: MB370_
Matrix (soil/w			_		le ID: MB370006
Level (low/med	·			_	eived: 07/13/96
		_		Date Reci	sived: 07/13/90
% Solids:		0			
Co	ncentration	Units (ug	/L or mg/kg dry	y weight)	: UG/L_
	CAS No.	Analyte	Concentration	C Q	М
	7440-38-2	Arsenic_	4.0		F_P_
	1	Cadmium_	3.3		P
	7440-47-3 7439-92-1	Chromium_ Lead	8.9	B	P_
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Color Before:	CLEAR	Clari	ty Before: CLE	AR_	Texture: N/A
Color After:	CLEAR	Clari	ty After: CLE	AR_	Artifacts:
Comments:					

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## GENERAL CHEMISTRY

#### CASE NARRATIVE GENERAL CHEMISTRY

QAL	Lab	Reference	No./SDG.	MB370

Project: Brown & Root Coastal Systems Station

#### I. RECEIPT

No exceptions were encountered unless a Sample Receipt Exception Report is attached to the Chain-of-Custody included with this data package.

#### II. HOLDING TIMES

All holding times were met.

#### III. METHOD

The method used is cited in the corresponding Form I.

#### IV. PREPARATION

Sample preparation proceeded normally, if applicable.

#### V. ANALYSIS

- A. Calibration: All acceptance criteria were met.
- B. Blanks: All acceptance criteria were met.
- C. Spikes: All acceptance criteria were met.
- D. Duplicates: All acceptance criteria were met.
- E. Laboratory Control Samples: All acceptance criteria were met.
- F. Samples: Sample analyses proceeded normally.

I certify that this data package is in compliance with the terms and conditions agreed to by the client and QAL, Inc., both technically and for completeness, except for the conditions noted above. Release of the data contained in this hardcopy data package has been authorized by the Laboratory Manager or designated person, as verified by the following signature.

SIGNED:

Velinda Herbert

General Organic/Inorganic Chemist

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Client Sample ID: 333-MW01-1 Sample Description: GRAB Sample Matrix: Water Date Collected: 07/41/96 (Thursday)
Date Received: 07/13/96 (Saturday)

CATEGORY NAME Analytical Parameter	Result	Units	Reporting Level	Date of Analysis	Analytical Method(s)
DEMAND AND GENERAL ORGANIC Total Petroleum Hydrocarbons	0.08	mg/L	0.05	07/23/96	EPA418.1
			: - <del></del>	,	a A

Client Sample ID: 333-MW02-1 Sample Description: GRAB Sample Matrix: Water

Date Collected: 07/11/96 (Thursday)
Date Received: 07/13/96 (Saturday)

C Analy	ATEGOR	Y NA	ME meter				Resu	ılt	Units	Rep	orting evel	Date of Analys	of sis	Analyt Method	ical (s)		<u> </u>
D Total	EMAND Petro	AND	GENERA Hydro	L ORGANI carbons	en en en en en en en en en en en en en e		1.1		mg/L		0.05	07/23/	196	EPA418	.1		
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Client Sample ID: 333-MW03-1 Sample Description: GRAB

Sample Matrix: Water

Date Collected: 07/11/96 (Thursday)
Date Received: 07/13/96 (Saturday)

CATEGORY NAME Analytical Parameter	Result	Units	Reporting Level	Date of Analysis	Analytical Method(s)
DEMAND AND GENERAL ORGANIC Total Petroleum Hydrocarbons	< 0.05	mg/L	0.05	07/23/96	EPA418.1
			(1) 作・数据 (1) た・1 は (1) よ・1	ME TO THE STATE OF	
	. 1	f	<del></del>	<del></del>	Vacation of the same of the sa

Client Sample ID: 333-PC4-1 Sample Description: GRAB Sample Matrix: Water Date Collected: 07/41/96 (Thursday)
Date Received: 07/13/96 (Saturday)

Lab Reference No: MB370 Lab Sample ID: MB370004

CATEGORY NAME Analytical Parameter	Result	Units	Reporting Date of Level Analysis	Analytical Method(s)
DEMAND AND GENERAL ORGANIC Total Petroleum Hydrocarbons	< 0.05	mg/L	0.05 07/23/96	EPA418.1
	4 40 4 10 4 	1 4 1 18	turing North Control	(6545

₫,ŧ

Client Sample ID: 333-MW01-1B Sample Description: GRAB Sample Matrix: Water

Date Collected: 07/11/96 (Thursday)
Date Received: 07/13/96 (Saturday)

CATEGORY NAME Analytical Parameter	Result	Units	Reporting Level	Date of Analysis	Analytical Method(s)
DEMAND AND GENERAL ORGANIC Total Petroleum Hydrocarbons	< 0.05	mg/L	0.05	07/23/96	EPA418.1

Client Sample ID: 333-MW03-1D Sample Description: GRAB Sample Matrix: Water

Date Collected: 07/11/96 (Thursday)
Date Received: 07/13/96 (Saturday)

CATEGORY NAME Analytical Parameter	Result	Units	Reporting Level	Date of Analysis	Analytical Method(s)
DEMAND AND GENERAL ORGANIC Total Petroleum Hydrocarbons	< 0.05	mg/L	0.05	07/23/96	EPA418.1
					÷
					1
	5 378gt, 10 2 <u>.</u>		Rest Lit		



## Brown & Root **Environmental**

455 FAIRWAY DRIVE, SUITE 200 DEERFIELD BEACH, FLORIDA 33441 (305) 570-5885 (305) 570-5974 (FAX)

SITE MANAGER: 6. 60000	SHIPPED TO: FAGE LOF_
PROJECT NAME: CTD 0008	apacity Analytical
BRE PROJECT NO.: 7113 CODE:	Montamery ALA
P.O. NO.:	(LABORATORY NAME, CITY)

CHAIN OF CUSTODY RECORD							L		OR											٦
			harlie Burgin	SAM	PΕ	RIX	PRES TYPE		1/4 1/4	- /	- /	- /	/	- /		\$ \&\$ / K	24 HR. [	]48 HR. [	TAT □ RUSH 172 HR. □ 7 DAY	
LAB NO.	DATE	TIME	SAMPLE IDENTIFICATION	COMP.	GRAB	MA	PARAMETERS		) 9/ v						2/0	RES		E DATE:		
	7-11-96	1714	333-GW-MWO1-001	¥	_	SW		3		2	$\perp$	1	2	1		SAU	ples =	PCEI	)	]
 		1602	333-6W-MW02-001		_	GN	11	Ц		Ш		$\perp \! \! \perp$	Ц	$\perp$						_
		1801	333-GW-MW03-001			54	Ш	Ц	Ш					$\perp$						4
		1530	333-PC4-155-001		_	GW	Ц	П		Ш		$\perp \parallel$	Д	Ш		FC	RLAE	B USE	ONLY	
<del></del>		1445	333-64-mwo1-001B		_	AFW	Ш	Ц				Ш	Ц			148# /	1B 3	70		
	1	1801	333-GW-MW03-001D		/	GW	1	1	٢.	+	4		L			PROJ#				
		1	Trip Blank													ACK				
			Temp Blank													HAZWRA CC LEVE		YN		
																CCEVE	LIZU	ICE		Π
																ANA REC		TEMP PH		П
																SAMPLE				7
			TOTAL NUMBER OF	CONT	ΓΑΙ	NERS														٦
1) Mickey	RELINQUISI	HED BY (SI	CNATURE) SEAL INTACT? DATE: 6/3	4/9	6	EMF	TY BO	TLES	RECI	EIVED	BY (	SIGN	ATURI	<u> </u>		SEA	L INTACT?	DAT	7-10-96	1
1) Whichen	( d).	Shine	(ES) NO N/A TIME: 9:	00A	M	(2)	EIVED	1	5					·				I/A TIME		
RELINQUISHED BY	Y (SIGNATI	JRE)	SEAL INTACT? DATE: 7-12  TES NO N/A TIME: 18	<u>-96</u>		(4)	_		SIGNA E	_	)						L'INTACT? NO N	DATI		-
RELINOUISHED BY	Y (SIGNATI	JRE)	SEAL INTACT? DATE:	5()_		REC	EIVEO	BY (	SIGNÁ	TURE'	)					SEA	L INTACT?	DATI	7-13-96	1
(5) YES NO N/A TIME:						6	On	ril		ير	<b>e</b>					(FES)	NO N	I/A TIME	09:00 AM	1
SPECIAL INSTRUCTIONS:						LAB	ORA	rof	₹Y F	REM	ARI	(S:								7
SAMPLE CONTAINERS PRECLEANED BY:  BRE LABORATORY EMANUFACTURER METHOD OF SHIPMEN						ed .	<b>%</b>					В	ILL	OF	LAD	ING NO.: _				]
WHITE-FULLY EXE YELLOW-RECEIVIN	CUTED CO	PY	SAMPLING JEAM:				RECE	IVED	FOR ATURE	LABOF	RATOR	lY J			•			No.	0063	1
PINK-SAMPLERS' GOLDENROD-SITE	COPY/QA	COPY				· · · · ·			7/13			<u>, 20c.</u>			c	9.00			_	

		Sample Receipt Exceptions Report
Batch	Number: MB 370	Origination date: 7/13/96
ال <sup>ة</sup> . ـ ـ ـ	avProject: Brown + Root 1	Environmental
SUN	MMARY OF EXCEPTION (chec	ck one if it applies)
1	Description of exception	Comments (write number of exception description and the impacted sample numbers)
	No custody seal as required by the project.	Sample ID 333-GW-PCY-155-001
	2. No chain-of-custody provided.	1 M 1 m 1 m 1 m 1 m 1 m 1 m 1 m 1 m 1 m
•	Chain-of-custody provided but incomplete.	one 504 EDB container is missing.
V	4. Samples broken or leaking on receipt.	Sample ID 333-6W-PCY-155-001
	5. Temperature of samples inappropriate for analysis requested.	lists two containers for 610 analysis,
	6. Container inappropriate for analysis requested.	but one container is broken.
	7. Inadequate sample volume for analysis requested.	
	<ol> <li>Preservation inappropriate for analysis requested.</li> </ol>	is .
	<ol> <li>Samples received out of holding time for analysis requested.</li> </ol>	
	10. Samples received more than 72 hours after sampling.	
	11. Discrepancies between chain-of- custody and container labels.	
V	12. Other (describe on right)	
70	ACTION(S) AFFECTED (*****	which fraction was affected by the exceptions detailed above by writing the number of the exception next to it)
	npreserved Nutric	
	Cyanide Extracta	ables Extractables Other (specify)
AC	CTION TAKEN:	

Cyanide	Extractables	Extractables	Other (specify)	·.
ACTION TAKEN:				
	:			
				· · · · · ·
		:		
Originator:		Supervisor:		
Client was notified on:	7/15	Client contact: Jerry	· boocle	
Client's comments:	Proceed of analy	us - add Pb from	e to metals	
		<del></del>		
		· · · · · · · · · · · · · · · · · · ·		
Client Con Jane		OA «Man-		<u></u>
Client Services:	Of m	QA officer:	γ <sup>8</sup> τ μ 3	

		Sampl	e Receipt E	exceptions	Report
Batch Number: MB370		Origination	n date: 7/15	196	
lent/Project: Room and Room	ŧ.	•			
SUMMARY OF EXCEPTIO	N (check one if	it applies)	-		
√ Description of exception		Comments (write number of excep	xion description and the	impacted sample nur	nibers)
1. No custody seal as required b	y the	. C. 1 11-41	-1.1-	-	•
project. 2. No classin-of-custody provided	i	ra Containers wout !	a bels.		
	7	sets of voa's		··	
3. Chain-of-custody provided bu incomplete.	5	EBB bottles		:	·
4. Samples broken or leaking or	receipt.	2 liter (1 baker)			
5. Temperature of samples inap for analysis requested.	propriate	metals ( liter) =		•	
6. Container inappropriate for a requested.	nalysis 3 -	- Ulter anhere (C10	)		
7. Inadequate sample volume for requested.	r analysis 2 -	1 liter aubers wide			
8. Preservation inappropriate for requested.	r analysis				
<ol> <li>Samples received out of hold for analysis requested.</li> </ol>	ing time				
10. Samples received more than after sampling.	72 hours	-	:		
11. Discrepancies between chair custody and container labels		;			
12. Other (describe on right)					
FRACTION(S) AFFECTED	(specify which fraction	was affected by the exceptions detaile	d above by writing the n	umber of the excention	on next to it)
Unpreserved	Nutrients	Metals		Volatiles	
Cyanide	Extractables	Extractables		Other (specify)	
ACTION TAKEN:					
	:	•	<del></del>	<del>,·_ ,</del>	
***************************************					
		•			

Originator:

Client was notified on: 7/15

Client contact: Terry Greek

Client's comments:

Client's comments:

Client Services:

QA officer:

## GROUNDWATER SAMPLES COLLECTED NOVEMBER 25, 1996

## CATIONS DATA PACKAGE

## CASE NARRATIVE Cations

Labor	atory:	CH2M HILL	Lab	Ref.	No.:	MC484					
Clien	t/Proje	ect:BROWN & ROOT CTO00008			<u></u>						
I.	<pre>Holding Time: All holding times were met.</pre>										
II.	<u>Digestion Exceptions</u> : None.										
III,	Analy	sis:									
	A.	<pre>Calibration: All acceptance criteria were met.</pre>									
	В.	Blanks: All acceptance criteria were met.									
	C.	<pre>ICP Interference Check Sample: All acceptance criteria were met.</pre>									
	D.	<pre>Spike Sample(s): All acceptance criteria were met.</pre>									
	E.	<pre>Duplicate Sample(s): All acceptance criteria were met.</pre>									
	F.	<u>Laboratory Control Sample(s)</u> : All acceptance criteria were met.									
	G.	<pre>ICP Serial Dilution: All acceptance criteria were met.</pre>									
	н.	Other: None.									
IV.	Any r	pt Exceptions: eceipt exception will be addressed t which will be attached to the Ch			_						
v.	Docum None.	entation Exceptions:				·					
VI.	condi Inc., detai packa	tify that this data package is in tions agreed to by the client and both technically and for complete led above. Release of the data coge has been authorized by the Laborified by the following signature.  D: Kaye Walker / Inorganic Division	Qual ness ntai rato	ity A , exc ned i ry Ma _ DAT	nalyt ept f n thi nager	ical Laboratories, for the conditions s hardcopy data					

EPA SAMPLE NO.

		INORGANIC A	ANALYSES DATA S	SHEET	1
مد. b Name: CH2M	HILL_		Contract: MC	C484	MW02-002
Lab Code: MGM_	Cas	se No.: MC4	184_ SAS No.:	: MC484_	SDG No.: MC484_
Matrix (soil/w	ater): WATE	R		Lab Sampl	e ID: MC484001
Level (low/med	): LOW	_		Date Rece	eived: 11/27/96
% Solids:	0.	0			
Co	ncentration	Units (ug	/L or mg/kg dry	y weight):	UG/L_
,	CAS No.	Analyte	Concentration	C Q	M
	7440-38-2 7440-43-9 7440-47-3 7439-92-1	CadmiumChromium	677		NR NR NR F
Calam Dafaara	D.D.OWN			_	
	BROWN		ty Before: CLE	_	Texture: N/A
	BROWN	Clari	ty After: CLE	AR_	Artifacts:
Comments:					

## 1

EPA SAMPLE NO

	:	INORGANIC A	ANALYSES DATA S	SHEET	1
.b Name: CH2M	HILL		Contract: MC	C484	TW01-001B
					SDG No.: MC484_
Matrix (soil/wa					le ID: MC484002
Level (low/med	): LOW_	_		Date Rec	eived: 11/27/96
% Solids:	0.	0			
Co	ncentration	Units (ug	/L or mg/kg dry	y weight)	: UG/L_
	1				
	CAS No.	Analyte	Concentration	C Q	M
•	7440-38-2	Arsenic_	1.6		F_
	7440-43-9		3.3		F_ P_ P_ F_
	7440-47-3				·   P
	7439-92-1	Lead	1.4	<sup>B</sup>	·  <sup>-</sup> -
	·			<del>-</del>	·   <del>  </del>
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				-	-
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	<del></del>				-
	<del></del>			-	-
				-	-
				-	-
Color Before:	CLEAR	Clari	ty Before: CLE	 AR	Texture: N/A
Color After:	CLEAR	Clari	ty After: CLE	AR_	Artifacts:
Comments:					

FORM I - IN

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EPA SAMPLE NO.

	-	ETIORGEMIT P	mmilloud Dain C	71111111	
.o Name: CH2M	HILL		Contract: MC		TW01-001
					SDG No.: MC484_
Matrix (soil/w					le ID: MC484003
Level (low/med	): LOW	_		Date Rece	eived: 11/27/96
% Solids:	0.0	0			
Co.	ncentration	Units (ug,	L or mg/kg dry	y weight):	: UG/L_
	CAS No.	Analyte	Concentration	C Q	М
	7440-47-3	Arsenic_Cadmium_Chromium_Lead	1.8 3.3 2.2 4.5	U	F P P F F F F F F F F F F F F F F F F F
Color Before:	YELLOW	Clari	lty Before: CLE	_  AR	Texture: N/A
Color After:			-	<del>-</del>	Artifacts:
Comments:					

### I INORGANIC ANALYSES DATA SHEET

EPA SAMPLE NO.

_ab Name: CH2M	_HILL		Contract: MC	C48	4	MW04-001
Lab Code: MGM_	Cas	se No.: MC4	84_ SAS No.:	: M	C484_	SDG No.: MC484_
Matrix (soil/w	ater): WATE	ર		La	b Sampl	e ID: MC484004
Level (low/med	): LOW	_		Da	te Rece	eived: 11/27/96
% Solids:	0.0	0				
Со	ncentration	Units (ug/	L or mg/kg dry	y w	eight):	UG/L_
	CAS No.	Analyte	Concentration	С	Q	М
	7440-47-3	Arsenic_Cadmium_Chromium_Lead	1.6 3.3 3.0 20.6	U B		F_P_P_F_
Color Before:	YELLOW	Clari	ty Before: CLO	UDY	?	Texture: N/A
Color After:	YELLOW	Clari	ty After: CLE	AR_	-	Artifacts:
Comments:						

## GENERAL CHEMISTRY

#### CASE NARRATIVE GENERAL CHEMISTRY

QAL Lab Reference No./SDG. MC484

Project: Brown & Root Coastal Systems Station

#### I. RECEIPT

No exceptions were encountered unless a Sample Receipt Exception Report is attached to the Chain-of-Custody included with this data package.

#### II. HOLDING TIMES

All holding times were met.

#### III. METHOD

The method used is cited in the corresponding Form I.

#### IV. PREPARATION

Sample preparation proceeded normally, if applicable.

#### V. ANALYSIS

- A. Calibration: All acceptance criteria were met.
- B. Blanks: All acceptance criteria were met.
- C. Spikes: There is no matrix spike data available because spiked sample was overrange and inadvertantly discarded. However, a post extraction matrix spike was performed and is included in this package.
- D. Duplicates: All acceptance criteria were met.
- E. Laboratory Control Samples: All acceptance criteria were met.
- F. Samples: Sample analyses proceeded normally.

I certify that this data package is in compliance with the terms and conditions agreed to by the client and QAL, Inc., both technically and for completeness, except for the conditions noted above. Release of the data contained in this hardcopy data package has been authorized by the Laboratory Manager or designated person, as verified by the following signature.

SIGNED:

Velinda Herbert

General Organic/Inorganic Chemist

å. å.

Client Sample ID: TW01-001B Sample Description: 333-GW Sample Matrix: Water Site: N/A

Date Collected: 11/25/96 12:15 (Mon)
Date Received: 11/27/96 10:00 (Wed)

Reference No: MC484 Lab Sample ID: MC484002

C. Analy	ATEGORY NA	AME ameter	-	Result	Units	Reporting Level	Date/Time of Analysis	Analytical Method(s)
D Total	EMAND AND Petroleum	GENERAL m Hydroca	ORGANIC arbons	< 0.06	mg/L	0.05	12/11/96 00:00	D EPA418.1
.**		-						
2 D								
				·			······································	(8196

W (8196)

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#### Report of Analytical Results

Client Sample ID: TW01-001 Sample Description: 333-GW Sample Matrix: Water Site: N/A Date Collected: 11/25/96 13:20 (Mon)
Date Received: 11/27/96 10:00 (Wed)

Reference No: MC484 Lab Sample ID: MC484003

CATEGORY NAME Analytical Parameter	Result	Units	Reporting Level	Date/Time of Analysis	Analytical Method(s)
DEMAND AND GENERAL ORGANI Total Petroleum Hydrocarbons	C 19.2	mg/L	0.25	12/11/96 00:00	EPA418.1
	ekut 1861 lunuk - Latin bilang La	an garagan kenalagan			The Shimpelia views (Cili
			d States in		
			환경성) - : 1 19 11		
· 			ing still The still		
	rinna (ing sa sa sa sa sa sa sa sa sa sa sa sa sa	of Medifical Leville House	arte un la fur		n na mara i Markeyanya kusake masa ne
				La Milia	
					4 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1

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#### Report of Analytical Results

Client Sample ID: MW04-001 Sample Description: 333-GW Sample Matrix: Water Site: N/A

Date Collected: 11/25/96 14:45 (Mon)
Date Received: 11/27/96 10:00 (Wed)

Reference No: MC484 Lab Sample ID: MC484004

CATEGORY NAME Analytical Parameter	Result	Units	Reporting Date/Time Level of Analysis	Analytical Method(s)
DEMAND AND GENERAL ORGANIC Total Petroleum Hydrocarbons	21.5	mg/L	0.51 12/11/96 00:	00 EPA418.1
			(2) (4) (2) (2) (2) (2) (3) (4) (4) (4) (4) (4) (4) (4) (4) (4) (4	

Ş .;

## CASE NARRATIVE GC/MS VOLATILE ORGANICS

QAL Lab	Reference	No./SDG.	MCMC	484	
Project:	Bro	wn & Root	<u>Coastal</u>	Systems	Station

#### I. RECEIPT

No exceptions were encountered unless a Sample Receipt Exception Report is attached to the Chain-of-Custody included with this data package.

#### II. HOLDING TIMES

- A. Sample Preparation: Not applicable.
- B. Sample Analysis: All holding times were met.

#### III. METHOD

Preparation: N/A Cleanup: N/A

Analysis: SW-846 8260

#### IV. PREPARATION

Not applicable.

#### V. ANALYSIS

- A. Calibration: All acceptance criteria were met.
- B. Blanks: All acceptance criteria were met.
- C. Surrogates: All acceptance criteria were met.
- D. Spikes: As requested, the matrix spikes were performed using a sample from sample delivery group MC484 (MC484003MS and MC484003MSD). All acceptance criteria were met. A copy of the results has been included for your review.
- E. Samples: Sample analysis proceeded normally.
- F. Other: Please note that the Form 1's reflect the specified target list.

Because this laboratory has only recently begun analyzing samples by the 8260 method, there are not yet enough data collected to produce control charts for the water surrogate recoveries for 1,2-Dichloroethane-d4. This control chart is currently in the process of being developed.

COLIMONO CHILLANDO CHOMILOS	
Lab Reference No./SDG.	MC484
Page 2	

I certify that this data package is in compliance with the terms and conditions agreed to by the client and QAL, Inc., both technically and for completeness except for the conditions noted above. Release of the data contained in this hardcopy data package has been authorized by the Laboratory Manager or designated person, as verified by the following signature.

SIGNED:

DATE:

Ward Dickens

Laboratory Director

Lab Reference No./SDG. MC484
Page 3

# CASE NARRATIVE Addendum

Sample Information

LAB	CLIENT	SAMPLE	DATE	DATE	DATE	SAMPLE
SAMPLE ID	SAMPLE ID	MATRIX	SAMPLED	EXTRACTED	ANALYZED	<sup>1</sup> Hq
MC484002	TW01-001B	WATER	11/25/96	N/A	12/04/96	<2
MC484003	TW01-001	WATER	11/25/96	N/A	12/04/96	<2
MC484003MS	TW01-001MS	WATER	11/25/96	N/A	12/04/96	<2
MC484003MSD	TW01-001MSD	WATER	11/25/96	N/A	12/04/96	<2
MC484004	MW04-001	WATER	11/25/96	N/A	12/06/96	<2
X12046B1	VBLKB3	WATER	N/A	N/A	12/04/96	N/A
X12066B1	VBLKB8	WATER	N/A	N/A	12/06/96	N/A

<sup>&</sup>lt;sup>1</sup> Applies to samples designated for purgeable VOA analysis only.

TW01-001B

Lab Name: CH2M HILL

Contract: MC484

Lab Code: MGM

Case No.: MC484

SAS No.: SDG No.: MC484

Matrix: (soil/water) WATER

Lab Sample ID: MC484002

Sample wt/vol: 5.0 (g/mL) ML

Lab File ID: C1VO040841.D

Level: (low/med) LOW

Date Received: 11/27/96

% Moisture: not dec. \_\_\_\_\_

Date Analyzed: 12/04/96

Column: (pack/cap) CAP

Dilution Factor: 1.0

CONCENTRATION UNITS:

(119/1, or 119/Kg) 119/1. CAS NO COMPOITNID

74-87-3	CAS NO.	COMPOUND (ug/L or t	ug/Kg) UG/L	Q
74-83-9Bromomethane       10       U         75-00-3				_
75-00-3Chloroethane       10       U         75-69-4Trichlorofluoromethane       10       U         75-35-41,1-Dichloroethene       10       U         75-09-2Methylene chloride       10       U         75-34-31,1-Dichloroethane       10       U         67-66-3Chloroform       10       U         71-55-6Chloroform       10       U         71-43-2Benzene       10       U         107-06-21,2-Dichloroethane       10       U         79-01-6Trichloroethene       10       U         78-87-51,2-Dichloropropane       10       U         75-27-4Bromodichloromethane       10       U         10061-01-5	74-83-9	Bromomethane	I	_
75-69-4Trichlorofluoromethane       10       U         75-35-41,1-Dichloroethene       10       U         75-09-2Methylene chloride       10       U         75-34-31,1-Dichloroethane       10       U         67-66-3Chloroform       10       U         71-55-61,1,1-Trichloroethane       10       U         56-23-5Carbon tetrachloride       10       U         71-43-2Benzene       10       U         107-06-21,2-Dichloroethane       10       U         79-01-6Trichloroethene       10       U         78-87-51,2-Dichloropropane       10       U         75-27-4Bromodichloromethane       10       U         100-75-82-Chloroethylvinyl ether       10       U         100-61-01-5cis-1,3-Dichloropropene       10       U         108-88-3Toluene       10       U         100-61-02-6trans-1,3-Dichloropropene       10       U         79-00-51,1,2-Trichloroethane       10       U         12-48-1Dibromochloromethane       10       U         108-90-7Chlorobenzene       10       U         100-41-4Bromoform       10       U <td>75-00-3</td> <td>Chloroethane</td> <td> I</td> <td></td>	75-00-3	Chloroethane	I	
75-35-41,1-Dichloroethene       10       U         75-09-2Methylene chloride       10       U         75-34-31,1-Dichloroethane       10       U         67-66-3Chloroform       10       U         71-55-61,1,1-Trichloroethane       10       U         56-23-5Carbon tetrachloride       10       U         71-43-2Benzene       10       U         107-06-21,2-Dichloroethane       10       U         79-01-6Trichloroethene       10       U         78-87-51,2-Dichloropropane       10       U         75-27-4Bromodichloromethane       10       U         100-61-01-5				_
75-09-2	75-35-4	1.1-Dichloroethene		_
75-34-31,1-Dichloroethane       10       U         67-66-3Chloroform       10       U         71-55-61,1,1-Trichloroethane       10       U         56-23-5Carbon tetrachloride       10       U         71-43-2Benzene       10       U         107-06-21,2-Dichloroethane       10       U         79-01-6Trichloroethene       10       U         78-87-51,2-Dichloropropane       10       U         75-27-4Bromodichloromethane       10       U         100-10-75-82-Chloroethylvinyl ether       10       U         108-88-3Toluene       10       U         108-88-3Toluene       10       U         100-10-2-6trans-1,3-Dichloropropene       10       U         79-00-51,1,2-Trichloroethane       10       U         127-18-4Tetrachloroethene       10       U         124-48-1Dibromochloromethane       10       U         108-90-7	75-09-2	Methylene chloride		
67-66-3Chloroform       10       U         71-55-61,1,1-Trichloroethane       10       U         56-23-5Carbon tetrachloride       10       U         71-43-2Benzene       10       U         107-06-21,2-Dichloroethane       10       U         79-01-6Trichloroethene       10       U         78-87-51,2-Dichloropropane       10       U         75-27-4Bromodichloromethane       10       U         110-75-82-Chloroethylvinyl ether       10       U         10061-01-5cis-1,3-Dichloropropene       10       U         108-88-3Toluene       10       U         10061-02-6trans-1,3-Dichloropropene       10       U         79-00-51,1,2-Trichloroethane       10       U         127-18-4Tetrachloroethene       10       U         124-48-1Dibromochloromethane       10       U         108-90-7	75-34-3	1,1-Dichloroethane		Ū
56-23-5Carbon tetrachloride       10       U         71-43-2	67-66-3	Chloroform	10	U.
71-43-2			10	IJ
107-06-21,2-Dichloroethane       10       U         79-01-6Trichloroethene       10       U         78-87-51,2-Dichloropropane       10       U         75-27-4Bromodichloromethane       10       U         110-75-82-Chloroethylvinyl ether       10       U         10061-01-5cis-1,3-Dichloropropene       10       U         108-88-3Toluene       10       U         109-00-51,1,2-Trichloroethane       10       U         127-18-4Tetrachloroethane       10       U         124-48-1Dibromochloromethane       10       U         108-90-7Chlorobenzene       10       U         100-41-4Ethylbenzene       10       U         75-25-2	56-23-5	Carbon tetrachloride	10	U
79-01-6Trichloroethene       10       U         78-87-51,2-Dichloropropane       10       U         75-27-4Bromodichloromethane       10       U         110-75-82-Chloroethylvinyl ether       10       U         10061-01-5cis-1,3-Dichloropropene       10       U         108-88-3Toluene       10       U         10061-02-6trans-1,3-Dichloropropene       10       U         79-00-51,1,2-Trichloroethane       10       U         127-18-4Tetrachloroethene       10       U         124-48-1Dibromochloromethane       10       U         108-90-7Chlorobenzene       10       U         100-41-4Ethylbenzene       10       U         75-25-2Bromoform       10       U			10	
78-87-51,2-Dichloropropane       10       U         75-27-4Bromodichloromethane       10       U         110-75-82-Chloroethylvinyl ether       10       U         10061-01-5cis-1,3-Dichloropropene       10       U         108-88-3Toluene       10       U         10061-02-6trans-1,3-Dichloropropene       10       U         79-00-51,1,2-Trichloroethane       10       U         127-18-4Tetrachloroethene       10       U         124-48-1Dibromochloromethane       10       U         108-90-7Chlorobenzene       10       U         100-41-4Ethylbenzene       10       U         75-25-2	107-06-2	1,2-Dichloroethane	10	
75-27-4Bromodichloromethane       10       U         110-75-82-Chloroethylvinyl ether       10       U         10061-01-5cis-1,3-Dichloropropene       10       U         108-88-3Toluene       10       U         10061-02-6trans-1,3-Dichloropropene       10       U         79-00-51,1,2-Trichloroethane       10       U         127-18-4Tetrachloroethene       10       U         124-48-1Dibromochloromethane       10       U         108-90-7Chlorobenzene       10       U         100-41-4Ethylbenzene       10       U         75-25-2Bromoform       10       U	79-01-6	Trichloroethene		
110-75-82-Chloroethylvinyl ether       10       U         10061-01-5cis-1,3-Dichloropropene       10       U         108-88-3Toluene       10       U         10061-02-6trans-1,3-Dichloropropene       10       U         79-00-51,1,2-Trichloroethane       10       U         127-18-4Tetrachloroethene       10       U         124-48-1Dibromochloromethane       10       U         108-90-7Chlorobenzene       10       U         100-41-4Ethylbenzene       10       U         75-25-2Bromoform       10       U	78-87-5	1,2-Dichloropropane	<del></del>	
10061-01-5cis-1,3-Dichloropropene       10       U         108-88-3Toluene       10       U         10061-02-6trans-1,3-Dichloropropene       10       U         79-00-51,1,2-Trichloroethane       10       U         127-18-4Tetrachloroethene       10       U         124-48-1Dibromochloromethane       10       U         108-90-7Chlorobenzene       10       U         100-41-4Ethylbenzene       10       U         75-25-2Bromoform       10       U	75-27-4	Bromodichloromethane		
108-88-3Toluene       10       U         10061-02-6trans-1,3-Dichloropropene       10       U         79-00-51,1,2-Trichloroethane       10       U         127-18-4Tetrachloroethene       10       U         124-48-1Dibromochloromethane       10       U         108-90-7Chlorobenzene       10       U         100-41-4Ethylbenzene       10       U         75-25-2Bromoform       10       U	110-75-8	2-Chloroethylvinyl ether	(	
10061-02-6trans-1,3-Dichloropropene       10       U         79-00-51,1,2-Trichloroethane       10       U         127-18-4Tetrachloroethene       10       U         124-48-1Dibromochloromethane       10       U         108-90-7Chlorobenzene       10       U         100-41-4Ethylbenzene       10       U         75-25-2Bromoform       10       U	10061-01-5	cis-1,3-Dichloropropene		
79-00-51,1,2-Trichloroethane       10       U         127-18-4Tetrachloroethene       10       U         124-48-1Dibromochloromethane       10       U         108-90-7Chlorobenzene       10       U         100-41-4Ethylbenzene       10       U         75-25-2Bromoform       10       U				
127-18-4Tetrachloroethene       10       U         124-48-1Dibromochloromethane       10       U         108-90-7Chlorobenzene       10       U         100-41-4Ethylbenzene       10       U         75-25-2Bromoform       10       U	10061-02-6	trans-1,3-Dichloropropene_	- <del></del> I	
124-48-1Dibromochloromethane       10       U         108-90-7Chlorobenzene       10       U         100-41-4Ethylbenzene       10       U         75-25-2Bromoform       10       U	79-00-5	1,1,2-Trichloroethane	1	
108-90-7Chlorobenzene       10       U         100-41-4Ethylbenzene       10       U         75-25-2Bromoform       10       U	127-18-4	Tetrachloroethene		
100-41-4Ethylbenzene 10 U 75-25-2Bromoform 10 U	124-48-1	Dibromochioromethane		
75-25-2Bromoform 10 U	108-90-7	Chloropenzene		
	100-41-4	Ecnylpenzene	<del></del>	
/9-34-51,1,2,2-letrachioroethane 10 0			I	~ ,
	/9-34-5	1,1,2,2-letrachioroethane_		U

### VOLATILE ORGANICS ANALYSIS DATA SHEET TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

TW01-001B

Lab Name: CH2M HILL

Contract: MC484

Lab Code: MGM

Case No.: MC484

SAS No.:

SDG No.: MC484

Matrix: (soil/water) WATER

Lab Sample ID: MC484002

Sample wt/vol: 5.0

(g/mL) ML

Lab File ID: C1VO040841.D

Date Received: 11/27/96

Level: (low/med)

LOW

% Moisture: not dec. \_\_\_

Data Analyzed: 12/04/96

Column: (pack/cap) CAP

Dilution Factor: 1.0

CONCENTRATION UNITS: Number TICs found: 0 (ug/L or ug/Kg) UG/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1. 2. 3.				
· · · - I				
6. —				
/ •				
9.				
10.				
12.				
14.				
10.				
16.				
19.				
20.				
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45. I				
27.				
29.				
		l		

#### 1A VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

TW01-001

Q

Lab Name: CH2M HILL Contract: MC484

Lab Code: MGM Case No.: MC484 SAS No.: SDG No.: MC484

Matrix: (soil/water) WATER Lab Sample ID: MC484003

Sample wt/vol: 5.0 (g/mL) ML Lab File ID: C2V0040842.D

Level: (low/med) LOW Date Received: 11/27/96

% Moisture: not dec. Date Analyzed: 12/04/96

Column: (pack/cap) CAP Dilution Factor: 1.0

CAS NO. COMPOUND CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L

74-87-3-----Chloromethane 10 U 75-01-4-----Vinyl chloride 10 U 74-83-9-----Bromomethane 10 U 75-00-3-----Chloroethane 10 U 75-69-4-----Trichlorofluoromethane 10 U 75-35-4-----1,1-Dichloroethene U 10 75-09-2-----Methylene chloride 10 U 75-34-3-----1,1-Dichloroethane 10 U U 67-66-3-----Chloroform 10 U 71-55-6-----1,1,1-Trichloroethane 10 U 56-23-5-----Carbon tetrachloride 10 J 71-43-2-----Benzene 2 107-06-2----1,2-Dichloroethane U 10 79-01-6-----Trichloroethene U 10 78-87-5----1,2-Dichloropropane 10 U 75-27-4-----Bromodichloromethane U 10 110-75-8----2-Chloroethylvinyl ether U 10 10061-01-5----cis-1,3-Dichloropropene U 10 108-88-3-----Toluene 10061-02-6-----trans-1,3-Dichloropropene U 10 10 U 79-00-5-----1,1,2-Trichloroethane 10 U 127-18-4-----Tetrachloroethene 10 U 124-48-1-----Dibromochloromethane 10 U U 108-90-7-----Chlorobenzene 10 100-41-4-----Ethylbenzene 10 U 75-25-2-----Bromoform 10 U 79-34-5-----1,1,2,2-Tetrachloroethane 10 U

### VOLATILE ORGANICS ANALYSIS DATA SHEET TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

TW01-001

Lab Name: CH2M HILL

Contract: MC484

Lab Code: MGM

Case No.: MC484 SAS No.:

SDG No.: MC484

Matrix: (soil/water) WATER

Lab Sample ID: MC484003

Sample wt/vol: 5.0

(q/mL) ML

Lab File ID:

C2VO040842.D

Date Received: 11/27/96

Level: (low/med)

% Moisture: not dec. \_\_\_\_

LOW

Data Analyzed: 12/04/96

Column: (pack/cap) CAP

Dilution Factor: 1.0

CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L

Number TICs found: 5

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1. 2. 3. 496-11-7 4. 5. 6. 7. 8. 9. 10.	Unknown Alkane Unknown Indane 1H-Indene dihydro methyl iso Unknown	17.067 18.667 18.850 21.050 21.817	7 9 18 10 7	===== J J
12. 13. 14. 15. 16. 17. 18.				
19. 20. 21. 22. 23. 24.				
26. 27. 28. 29.				

## VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

MW04-001

Lab Name: CH2M HILL

Contract: MC484

Lab Code: MGM Case No.: MC484 SAS No.:

SDG No.: MC484

Matrix: (soil/water) WATER

Lab Sample ID: MC484004

Sample wt/vol:

Lab File ID: C2VO040864.D

5.0 (g/mL) ML

Level: (low/med) LOW

Date Received: 11/27/96

% Moisture: not dec.

Date Analyzed: 12/06/96

Column: (pack/cap) CAP

Dilution Factor: 1.0

CONCENTRATION UNITS:

CAS NO.

COMPOUND

(ug/L or ug/Kg) UG/L

Q

		.,	
74-87-3	Chloromethane	10	ן ד
75-01-4	Vinyl chloride	10	U
74-83-9	Bromomethane	10	ט
75-00-3	Chloroethane	10	ן ני
75-69-4	Trichlorofluoromethane	10	ט
75-35-4	1,1-Dichloroethene	10	ט
75-09-2	Methylene chloride	10	ט
75-34-3	1,1-Dichloroethane	10	ט
	Chloroform	10	ט (
71-55-6	1,1,1-Trichloroethane	10	ט
56-23-5	Carbon tetrachloride	10	ט ו
71-43-2	Benzene	3	J
	1,2-Dichloroethane	10	ט ו
79-01-6	Trichloroethene	10	ע
78-87-5	1,2-Dichloropropane	10	ע
75-27-4	Bromodichloromethane	10	י
110-75-8	2-Chloroethylvinyl ether	10	ט ט
10061-01-5	cis-1,3-Dichloropropene	10:	U
108-88-3	Toluene	10	Ū
	trans-1,3-Dichloropropene	10	Ū
79-00-5	1,1,2-Trichloroethane	10	Ū
127-18-4	Tetrachloroethene	10	บ
	Dibromochloromethane	10	Ū
	Chlorobenzene	10	Ū
100-41-4	Ethylbenzene	7	J
75-25-2	Bromoform	10	บิ
	1,1,2,2-Tetrachloroethane	10	Ū
		· · · · · · · · · · · · · · · · · · ·	

#### 1E VOLATILE ORGANICS ANALYSIS DATA SHEET TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

MW04-001

Lab Name: CH2M HILL

Contract: MC484

Lab Code: MGM Case No.: MC484 SAS No.: SDG No.: MC484

Matrix: (soil/water) WATER

Lab Sample ID: MC484004

Sample wt/vol: 5.0 (g/mL) ML

Lab File ID: C2VO040864.D

Level: (low/med) LOW

Date Received: 11/27/96

% Moisture: not dec.

Data Analyzed: 12/06/96

Column: (pack/cap) CAP

Dilution Factor: 1.0

Number TICs found: 9

CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
			==========	
1.	Unknown Alkane	17.083	65	J
2.	Unknown Alkane	17.817	90	J
3.	Unknown	18.767	110	J
3.	Unknown Aromatic Hydrocarbon		48	J
5.	Unknown Alkane	19.033		
		19.633	52	J
6. 119-64-2	Naphthalene, 1,2,3,4-tetrahy	21.433	66	ИJ
7.	Naphthalene tetrahydro methy	23.183	61	J
8.	Unknown Aromatic Hydrocarbon		66	J
9.	Naphthalene methyl isomer	24.017	80	J
10			<u> </u>	
11.				
14				
1.J.				
15. 16. 17. 18.				
16.				
17.				
18.	· · · · · · · · · · · · · · · · · · ·			
19	<del></del>			
19.				
20.				
21.				
22.				
23.				
44.				
47.				
20.		i		
41.				
20.				
47.				
30				

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## SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: CH2M HILL

Contract: MC484

MW04-001

Lab Code: MGM

Case No.: MC484 SAS No.:

SDG No.: MC484

Matrix: (soil/water) WATER

Lab Sample ID: MC484004

Sample wt/vol: 1000 (g/mL) ML

Lab File ID:

05DEC0701007.D

Level: (low/med) LOW

Date Received: 11/27/96

% Moisture: not dec. \_\_\_\_\_ dec.\_\_\_

Date Extracted:11/30/96

Extraction: (SepF/Cont/Sonc) CONT

Date Analyzed: 12/05/96

GPC Cleanup: (Y/N) N

pH: 7.0

Dilution Factor: 10.0

CONCENTRATION UNITS:

CAS NO.

COMPOUND

(ug/L or ug/Kg) UG/L

Q

534-52-14,6-Dinitro-2-methylphenol	500	ט
86-30-6N-Nitrosodiphenylamine $(\overline{3})$	100	1
122-66-71,2-Diphenylhydrazine	100	
101-55-34-Bromophenyl-phenylether	100	
118-74-1Hexachlorobenzene	100	ט
87-86-5Pentachlorophenol	500	ט ו
85-01-8Phenanthrene	160	
120-12-7Anthracene	24	
84-74-2Di-n-butylphthalate	100	
206-44-0Fluoranthene	11	J
92-87-5Benzidine	100	ן די
129-00-0Pyrene	41	J
85-68-7Butylbenzylphthalate	100	U
56-55-3Benzo (a) anthracene	100	ָּט
91-94-13,3'-Dichlorobenzidine	200	U
218-01-9Chrysene	100	U
117-81-7bis(2-Ethylhexyl)phthalate	12	JB
117-84-0Di-n-octylphthalate	100	U
205-99-2Benzo (b) fluoranthene	100	ָּט
207-08-9Benzo(k) fluoranthene	100	[ ט
50-32-8Benzo (a) pyrene	100	Ū
193-39-5Indeno (1, 2, 3-cd) pyrene	100	Ū
53-70-3Dibenz (a, h) anthracene	100	IJ
191-24-2Benzo(g,h,i)perylene	100	ט

(3) - Cannot be separated from Diphenylamine

FORM I SV-2

SW846

MW04-001

Lab Name: CH2M HILL

Contract: MC484

Lab Code: MGM Case No.: MC484 SAS No.: SDG No.: MC484

Matrix: (soil/water) WATER

Lab Sample ID: MC484004

Sample wt/vol: 1000 (q/mL) ML Lab File ID: 05DEC0701007.D

Date Received: 11/27/96

Level: (low/med) LOW

% Moisture: not dec.\_\_\_\_ dec.\_\_\_

Date Extracted:11/30/96

Extraction: (SepF/Cont/Sonc) CONT

Date Analyzed: 12/05/96

GPC Cleanup: (Y/N) N pH: 7.0 Dilution Factor: 10.0

CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L

Number TICs found: 20

CAC ATROPED	GOMBOTHE MANGE	D.T.	TIGTH GOVIG	
CAS NUMBER	COMPOUND NAME	K.T.	EST. CONC.	Q
CAS NUMBER  ===================================	COMPOUND NAME  ===================================	7.474 7.667 7.845 8.160 8.345 8.431 8.667 9.260 9.432 10.432 10.539 10.596	190	ר כר כר כר כר כר מת מת בר כר בר בר בר בר בר בר בר בר בר בר בר בר בר
18. 19.	Unknown Alkane Unknown PAH	12.104 12.525	500 260	J J

GC EXTRACTABLE VOLATILE ORGANICS (EDB)

# CASE NARRATIVE GC EXTRACTABLE VOLATILE ORGANICS (EDB)

QAL Lab F	Reference	No./SDG.	MC484		•
Project:	Brown &	Root Coa	stal Sys	tems Statio	n

#### I. RECEIPT

No exceptions were encountered unless a Sample Receipt Exception Report is attached to the Chain-of-Custody included with this data package.

#### II. HOLDING TIMES

- A. Sample Preparation: All holding times were met.
- B. Sample Analysis: All holding times were met.

#### III. METHOD

Preparation: N/A Cleanup: N/A Analysis: EPA 504.1

#### IV. PREPARATION

Sample preparation proceeded normally.

#### V. ANALYSIS

- A. Calibration: All acceptance criteria were met.
- B. Blanks: All acceptance criteria were met.
- C. Surrogates: All acceptance criteria were met.
- D. Spikes: Matrix spikes were performed using a sample from this contract. The associated blank spike (LCS) is summarized with the MS/MSD results on the Form 3 in this report.
- E. Samples: Sample analysis proceeded normally.

I certify that this data package is in compliance with the terms and conditions agreed to by the client and QAL, Inc., both technically and for completeness except for the conditions noted above. Release of the data contained in this hardcopy data package has been authorized by the Laboratory Manager or designated person, as verified by the following signature.

SIGNED: Tamon Caren	DATE: 12/23/96
Tammy Carey	
Chemist	

4.

GC EXTRACTABLE VOLATILE ORGANICS (EDB)
Lab Reference No./SDG: MC484
Page 2

# CASE NARRATIVE Addendum

Sample Information

LAB SAMPLE ID	CLIENT SAMPLE ID	SAMPLE MATRIX	DATE SAMPLED	DATE EXTRACTED	DATE ANALYZED	SAMPLE pH <sup>1</sup>
MC484002	TW01-001B	WATER	11/25/96	12/20/96	12/20/96	N/A
MC484003	TW01-001	WATER	11/25/96	12/20/96	12/20/96	N/A
MC484003MS	TW01-001MS	WATER	11/25/96	12/20/96	12/20/96	N/A
MC484003MSD	TW01-001MSD	WATER	11/25/96	12/20/96	12/20/96	N/A
MC484004	MW04-001	WATER	11/25/96	12/20/96	12/20/96	N/A
W12206B1	QC BLANK	WATER	N/A	12/20/96	12/20/96	N/A

<sup>1</sup> Applies to samples designated for purgeable VOA analysis only.

Laboratory Name: CH2M HILL Concentration: LOW Date Extracted: 12/20/96
Lab Sample ID: MC484002 Sample Matrix: WATER Date Analyzed: 12/20/96
Client Sample ID: TW01-001B Percent Moisture: Dilution Factor: 1.0

EDB

CAS Number uq/L
106-93-4 1,2-Dibromoethane (EDB) . . 0.02 U

1,1,2,2-Tetrachloroethane - SS 96

- U Analyzed for but not detected.
- B Detected in QC blank.
- J Detected, concentration estimated.
- SS Surrogate Standard reported as percent recovery.

Comments:

Form I

\$ 1

TLC

Laboratory Name: CH2M HILL Concentration: LOW Date Extracted: 12/20/96
Lab Sample ID: MC484003 Sample Matrix: WATER Date Analyzed: 12/20/96
Client Sample ID: TW01-001 Percent Moisture: Dilution Factor: 1.0

EDB

CAS Number ug/L
106-93-4 1,2-Dibromoethane (EDB) . . . 0.02 U

1,1,2,2-Tetrachloroethane - SS 108

U - Analyzed for but not detected.

B - Detected in QC blank.

J - Detected, concentration estimated.

SS - Surrogate Standard reported as percent recovery.

Comments:

Form I

4. 4

Laboratory Name	: CH2M HILL	Concentration:	LOW	Date Extracted:	12/20/96
Lab Sample ID:	MC484004	Sample Matrix:	WATER_	Date Analyzed:	12/20/96
Client Sample I	D: MW04-001	Percent Moisture:		Dilution Factor:	1.0

EDB

<u>CAS Number</u> <u>uq/L</u> 106-93-4 1,2-Dibromoethane (EDB) . . . 0.02 U

. \*

1,1,2,2-Tetrachloroethane - SS 100

- U Analyzed for but not detected.
- B Detected in QC blank.
- J Detected, concentration estimated.
- SS Surrogate Standard reported as percent recovery.

Comments:

Form I

4, 4

GC POLYNUCLEAR AROMATIC HYDROCARBONS

## CASE NARRATIVE GC POLYNUCLEAR AROMATIC HYDROCARBONS

QAL Lab	Reference	No./SDG.	MC4	84	•	<u>-</u>	 	 
Project:	Brown &	Root Coas	tal	Svstems	Station			

#### I. RECEIPT

No exceptions were encountered unless a Sample Receipt Exception Report is attached to the Chain-of-Custody included with this data package.

#### II. HOLDING TIMES

- A. Sample Preparation: All holding times were met.
- B. Sample Analysis: All holding times were met.

#### III. METHOD

Preparation: N/A Cleanup: N/A Analysis: 610

#### IV. PREPARATION

Sample preparation proceeded normally.

#### V. ANALYSIS

A. Calibration: All acceptance criteria were met.

Both the initial calibration and continuing calibration summaries include data for both the primary and confirmation columns. Each compound will appear in the summary reports twice. The first time the compound will not be proceeded by the "#" symbol, referring to compounds identified from the first column (RTX-5); the next time it will have the "#"symbol, referring to compounds identified from the second column (RTX-200) (for example, Naphthalene and #Naphthalene).

- B. Blanks: All acceptance criteria were met.
- C. Surrogates: Surrogate recovery could not be determined for sample MC484004 due to the dilution required for analysis. All other acceptance criteria were met.
- D. Spikes: Matrix spikes were performed using a sample from this contract. MC484003MS and MC484003MSD were diluted for analysis due to native contents. Therefore some recoveries are inflated by coeluting non-target peaks. Recovery of Benzo(k)fluoranthene could not be determined due to the dilution required for analysis. The associated blank spike (LCS) is additionally summarized on the Form 3 in this report.

E. Samples: Samples MC484003, MC484004, MC484003MS, and MC484003MSD were diluted for analysis due to non-target interference. Sample analysis proceeded normally.

The report limit for Phenanthrene was raised for sample MC484004 due to chromatographic interference.

Frequently interferences will persist in the extract even after cleanup procedures. Standard cleanup procedures are designed to recover the target compounds and remove interfering non-targets. Precise rules for diluting interferences are difficult to develop. A single non-target peak could be allowed to saturate the detector. However, extracts with multiple non-target peaks might elevate the baseline or alter the baseline noise for part of the chromatogram. If the baseline were severely elevated or noise obstructed the target chromatographic region, targets at or near the report limit could not be positively identified within the interfering peaks. Because GC identification is based largely on retention time, regions with many peaks (noisy regions) will frequently have many false positives. For samples with such chromatographic interference, positive hits are typically not reported unless the peak is significantly above the surrounding noise and/or is not obstructed on one or both analytical columns. In addition, such interferences can be damaging to the chromatography and on-going calibration criteria can not be achieved. Without dilution, report limits would usually be raised in samples with significant interference. Therefore, such samples are typically diluted to minimize interferences and yet achieve the best possible report limits.

F. Other: Primary and confirmation data were simultaneously acquired using two dissimilar analytical columns (RTX-5 and RTX-200) connected in parallel to one injection port and one detector.

I certify that this data package is in compliance with the terms and conditions agreed to by the client and QAL, Inc., both technically and for completeness except for the conditions noted above. Release of the data contained in this hardcopy data package has been authorized by the Laboratory Manager or designated person, as verified by the following signature.

SIGNED: Jammy (A)
Tammy Carey

Chemist

DATE: 12/23/96

GC POLYNUCLEAR AROMATIC HYDROCARBONS Lab Reference No./SDG: MC484 Page

### CASE NARRATIVE Addendum

Sample Information

LAB SAMPLE ID	CLIENT SAMPLE ID	SAMPLE MATRIX	DATE SAMPLED	DATE EXTRACTED	DATE ANALYZED	SAMPLE pH <sup>1</sup>
MC484002	TW01-001B	WATER	11/25/96	11/30/96	12/18/96	N/A
MC484003	TW01-001	WATER	11/25/96	11/30/96	12/18/96	N/A
MC484003MS	TW01-001MS	WATER	11/25/96	11/30/96	12/18/96	N/A
MC484003MSD	TW01-001MSD	WATER	11/25/96	11/30/96	12/18/96	N/A
MC484004	MW04-001	WATER	11/25/96	11/30/96	12/18/96	N/A
W11306B1	QC BLANK	WATER	N/A	11/30/96	12/18/96	N/A

<sup>1</sup> Applies to samples designated for purgeable VOA analysis only.

Laboratory Name: CH2M HILL Concentration: LOW Date Extracted: 11/30/96
Lab Sample ID: MC484002 Sample Matrix: WATER Date Analyzed: 12/18/96
Client Sample ID: TW01-001B Volume Extracted: 1025mL Dilution Factor: \_\_\_\_\_1.0

#### PNA COMPOUNDS

CAS Number		uq/L	
91-20-3	Naphthalene	2	U
91-57-6	2-Methylnaphthalene	2	U
90-12-0	1-Methylnaphthalene	2	U
208-96-8	Acenaphthylene	2	U
83-32-9	Acenaphthene	2	U
86-73-7	Fluorene	2	U
85-01 <b>-</b> 8	Phenanthrene	2	U
120-12-7	Anthracene	2	υ
206-44-0	Fluoranthene	2	U
129-00-0	Pyrene	2	U
56-55-3	Benzo(a) anthracene	2	U
218-01-9	Chrysene	2	Ω
205-99-2	Benzo(b) fluoranthene	2	U
207-08-9	Benzo(k)fluoranthene	2	U
50-32-8	Benzo(a)pyrene	2	U
193-39-5	Indeno(1,2,3-cd)pyrene	2	U
53-70-3	Dibenzo(a,h)anthracene	2	U
191-24-2	Benzo(g,h,i)perylene	2	U
	Terphenyl-d14 - SS	66	ક

#### Comments:

Form I

U - Analyzed for but not detected.

B - Detected in QC blank.

J - Detected, concentration estimated.

SS - Surrogate Standard reported as percent recovery.

Laboratory Name: CH2M HILL Concentration: LOW Date Extracted: 11/30/96
Lab Sample ID: MC484003 Sample Matrix: WATER Date Analyzed: 12/18/96
Client Sample ID: TW01-001 Volume Extracted: 990mL Dilution Factor: \_\_\_\_5.0

#### PNA COMPOUNDS

CAS Number		uq/L	
91-20-3	Naphthalene	10	U
91-57-6	2-Methylnaphthalene	10	U
90-12-0	1-Methylnaphthalene	10	U
208-96-8	Acenaphthylene	10	υ
83-32-9	Acenaphthene	10	U
86-73-7	Fluorene	10	U
85-01-8	Phenanthrene	10	U
120-12-7	Anthracene	10	U
206-44-0	Fluoranthene	10	U
129-00-0	Pyrene	10	U
56-55-3	Benzo(a) anthracene	10	U
218-01-9	Chrysene	10	U
205-99-2	Benzo(b) fluoranthene	10	U
207-08-9	Benzo(k)fluoranthene	10	U
50-32-8	Benzo(a)pyrene	10	U
193-39-5	Indeno(1,2,3-cd)pyrene	10	U
53-70-3	Dibenzo(a,h)anthracene	10	U
191-24-2	Benzo(g,h,i)perylene	10	U
	Terphenyl-d14 - SS	60	%

#### Comments:

Form I

U - Analyzed for but not detected.

B - Detected in QC blank.

J - Detected, concentration estimated.

SS - Surrogate Standard reported as percent recovery.

Concentration: Laboratory Name: CH2M HILL LOW Date Extracted: Date Analyzed: Lab Sample ID: MC484004 Sample Matrix: WATER 12/18/96 Client Sample ID: MW04-001 Volume Extracted: 1031mL Dilution Factor:

#### PNA COMPOUNDS .

CAS Number		ug/L	
91-20-3	Naphthalene	200	U
91-57-6	2-Methylnaphthalene	340	
90-12-0	1-Methylnaphthalene	330	
208-96-8	Acenaphthylene	200	U
83-32-9	Acenaphthene	200	U
86-73-7	Fluorene	200	U
85-01-8	Phenanthrene	400	U.
120-12-7	Anthracene	200	U
206-44-0	Fluoranthene	200	U
129-00-0	Pyrene	200	U
56-55-3	Benzo(a) anthracene	200	U
218-01-9	Chrysene	200	υ
205-99-2	Benzo(b) fluoranthene	200	U
207-08-9	Benzo(k) fluoranthene	200	U
50-32-8	Benzo(a)pyrene	200	Ω
193-39-5	Indeno(1,2,3-cd)pyrene	200	U
53-70-3	Dibenzo(a,h)anthracene	200	Ū
191-24-2	Benzo(g,h,i)perylene	200	U
	Terphenyl-d14 - SS	DL	

Comments: DL - Surrogate recovery not determined due to dilution.

Form I

U - Analyzed for but not detected.

B - Detected in QC blank.

J - Detected, concentration estimated.

SS - Surrogate Standard reported as percent recovery.

Chain of custody documentation



### Brown & Root Environmental

455 FAIRWAY DRIVE, SUITE 200 DEERFIELD BEACH, FLORIDA 33441 (305) 570-5885 (305) 570-5974 (FAX)

SITE MANAGER: _	Gerald	F. 600	de	
PROJECT NAME:	CTO	0008	Coash	systems bhabie

BRE PROJECT NO.: \_\_\_\_\_ CODE: \_\_\_\_

P.O. NO.: 2049-7113-P96249

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Montgomery, AL

(LABORATORY NAME, CITY)

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QAL Montgomery

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PLERECT.DOC (V20%)

GC PURGEABLE HALOCARBONS

## CASE NARRATIVE GC PURGEABLE HALOCARBONS

QAL Lab	Reference	No./SDO	G: MC4	84			
Project:	Brown	& Root	Coastal	Systems	Station		

#### I. RECEIPT

No exceptions were encountered unless a Sample Receipt Exception Report is attached to the Chain-of-Custody included with this data package.

#### II. HOLDING TIMES

- A. Sample Preparation: N/A
- B. Sample Analysis: All holding times were met.

#### III. METHOD

. Preparation:

Cleanup: N/A

Analysis: EPA 601 (Mod)

#### IV. PREPARATION

Not applicable.

#### V. ANALYSIS

- A. Calibration: All acceptance criteria were met.
- B. Blanks: All acceptance criteria were met.
- C. Surrogates: All acceptance criteria were met.
- D. Spikes: Recovery of 2-chloroethyl vinyl ether (2-CEVE) was low in both matrix spikes which is not unusual for acid preserved samples. The laboratory control sample analyzed immediately after the matrix spikes was not preserved with acid and provided good recovery of 2-CEVE. All other recoveries were within the advisory limits.
- E. Samples: Sample analysis proceeded normally. Primary analysis was performed using a J&W DB-VRX analytical column (75 m x 0.45 mm ID). Confirmational analyses were performed using a Restek Rtx-502.2 analytical column (105 m x 0.53 mm ID).

I certify that this data package is in compliance with the terms and conditions agreed to by the client and QAL, Inc., both technically and for completeness except for the conditions noted above. Release of the data contained in this hardcopy data package has been authorized by the Laboratory Manager or designated person as verified by the following signature.

SIGNED: Mult W. Sywloy DATE: 12-20-96

Herb Kelly
Organic Division Manager

Page 2

### CASE NARRATIVE Addendum

Sample Information

LAB SAMPLE ID	CLIENT SAMPLE ID	SAMPLE MATRIX	DATE SAMPLED	DATE EXTRACTED	DATE ANALYZED	SAMPLE pH <sup>1</sup>
MC484002	TW01-001B	WATER	11/25/96	N/A	12/02/96	<2
MC484003	TW01-001	WATER	11/25/96	N/A	12/02/96	<2
MC484003MS	TW01-001MS	WATER	11/25/96	N/A	12/02/96	<2
MC484003MSD	TW01-001MSD	WATER	11/25/96	N/A	12/02/96	<2
MC484004	MW04-001	WATER	11/25/96	N/A	12/02/96	<2
X12026B1	VBLK001	WATER	N/A	N/A	12/02/96	N/A

 $<sup>^{\</sup>scriptsize 1}$  Applies to samples designated for purgeable VOA analysis only.

# CURRENT METHOD DETECGTION LIMITS (MDLs) PURGEABLE HALOCARBONS

Date Collected: N/A Sample Group: Lab QC

Date Extracted: N/A Lab Sample ID: Multiple Samples

Date Analyzed: 12/20/96 Lab File 1 ID: N/A
Matrix: Water Lab File 2 ID: N/A
Method: EPA 601 (Mod) Dilution Factor: 1.0

% Moisture: 100 Reporting Units: ug/L

CAS NUMBER	COMPOUND NAME	REPORTING LIMIT	RESULT
75-27-4	Bromodichloromethane	1.0	0.178
75-25-2	Bromoform	1.0	0.362
74-83-9	Bromomethane	1.0	0.143
56-23-5	Carbon tetrachloride	1.0	0.191
108-90-7	Chlorobenzene	1.0	0.746
75-00-3	Chloroethane	1.0	0.085
110-75-8	2-Chloroethyl vinyl ether	1.0	0.059
67-66-3	Chloroform	1.0	0.064
74-87-3	Chloromethane	1.0	0.148
124-48-1	Dibromochloromethane	1.0	0.069
95-50-1	1,2-Dichlorobenzene	1.0	0.298
541-73-1	1,3-Dichlorobenzene	1.0	0.319
106-46-7	1,4-Dichlorobenzene	1.0	0.358
75-71-8	Dichlorodifluoromethane	1.0	0.195
75-34-3	1,1-Dichloroethane	1.0	0.128
107-06-2	1,2-Dichloroethane	1.0	0.043
75-35-4	1,1-Dichloroethene	1.0	0.125
156 <b>-</b> 59-2	cis-1,2-Dichloroethene	1.0	0.054
156-60-5	trans-1,2-Dichloroethene	1.0	0.153
78-87-5	1,2-Dichloropropane	1.0	0.155
10061-01-5	cis-1,3-Dichloropropene	1.0	0.091
10061-02-6	trans-1,3-Dichloropropene	1.0	0.038
75-09-2	Methylene chloride	5.0	0.512
79-34-5	1,1,2,2-Tetrachlorethane	1.0	0.113
127-18-4	Tetrachloroethene	1.0	0.281
71-55-6	1,1,1-Trichloroethane	1.0	0.173
79-00-5	1,1,2-Trichloroethane	1.0	0.077
79-01-6	Trichloroethene	1.0	0.147
75-69-4	Trichlorofluoromethane	1.0	0.112
75-01-4	Vinyl chloride	1.0	0.180



# REPORT OF ANALYTICAL RESULTS PURGEABLE HALOCARBONS

Date Collected: 11/25/96

Date Extracted: N/A

Date Extracted: N/A

Date Analyzed: 12/02/96

Matrix: Water

Method: EPA 601 (Mod)

Sample Group: MC484

Lab Sample ID: MC484002

Lab File 1 ID: N02V012

Lab File 2 ID: N02W012

Dilution Factor: 1.0

% Moisture: 100 Reporting Units: ug/L

CAS NUMBER	COMPOUND NAME	REPORTING LIMIT	RESULT
75-27-4	Bromodichloromethane	1.0	U
75-25-2	Bromoform	1.0	ប
74-83-9	Bromomethane	1.0	ט
56-23-5	Carbon tetrachloride	1.0	ប
108-90-7	Chlorobenzene	1.0	ט
75-00-3	Chloroethane	1.0	ប
110-75-8	2-Chloroethyl vinyl ether	1.0	U
67-66-3	Chloroform	1.0	ប
74-87-3	Chloromethane	1.0	U
124-48-1	Dibromochloromethane	1.0	Ū
95-50-1	1,2-Dichlorobenzene	1.0	U
541-73-1	1,3-Dichlorobenzene	1.0	U
106-46-7	1,4-Dichlorobenzene	1.0	U
75-71-8	Dichlorodifluoromethane	1.0	U
75-34-3	1,1-Dichloroethane	1.0	ט
107-06-2	1,2-Dichloroethane	1.0	ប
75-35-4	1,1-Dichloroethene	1.0	U
156-59-2	cis-1,2-Dichloroethene	1.0	ប
156-60-5	trans-1,2-Dichloroethene	1.0	ប
78-87-5	1,2-Dichloropropane	1.0	ប
10061-01-5	cis-1,3-Dichloropropene	1.0	U
10061-02-6	trans-1,3-Dichloropropene	1.0	ប
75-09-2	Methylene chloride	5.0	ប
79-34-5	1,1,2,2-Tetrachlorethane	1.0	ប
127-18-4	Tetrachloroethene	1.0	ប
71-55-6	1,1,1-Trichloroethane	1.0	U
79-00-5	1,1,2-Trichloroethane	1.0	υ
79-01-6	Trichloroethene	1.0	U
75-69-4	Trichlorofluoromethane	1.0	ប
75-01-4	Vinyl chloride	1.0	υ
/3-01-4	vinyi chioride	1.0	

SURROGATE-Fluorobenzene (QC Limits - 61-133%)

101 % Rec.



# REPORT OF ANALYTICAL RESULTS PURGEABLE HALOCARBONS

Date Collected: 11/25/96

Date Extracted: N/A
Date Analyzed: 12/02/96

Matrix: Water

Method: EPA 601 (Mod) % Moisture: 100

Sample Group: MC484

Lab Sample ID: MC484003 Lab File 1 ID: N02V013 Lab File 2 ID: N02W013

Dilution Factor: 1.0

Reporting Units: ug/L

CAS NUMBER	COMPOUND NAME	REPORTING LIMIT	RESULT	
75-27-4	Bromodichloromethane	1.0	ŭ	
75-25-2	Bromoform	1.0	U	
74-83-9	Bromomethane	1.0	บ	
56-23-5	Carbon tetrachloride	1.0	U	
108-90-7	Chlorobenzene	1.0	บ	
75-00-3	Chloroethane	1.0	บ	
110-75-8	2-Chloroethyl vinyl ether	1.0	U	
67-66-3	Chloroform	1.0	U	
74-87-3	Chloromethane	1.0	ָ ט	
124-48-1	Dibromochloromethane	1.0	υ	
95-50-1	1,2-Dichlorobenzene	1.0	บ	
541-73-1	1,3-Dichlorobenzene	1.0	υ	
106-46-7	1,4-Dichlorobenzene	1.0	ט	
75-71-8	Dichlorodifluoromethane	1.0	υ	
75-34-3	1,1-Dichloroethane	1.0	υ	
107-06-2	1,2-Dichloroethane	1.0	บ	
75-35-4	1,1-Dichloroethene	1.0	ָּט	
156-59-2	cis-1,2-Dichloroethene	1.0	บ	
156-60-5	trans-1,2-Dichloroethene	1.0	υ	
78-87-5	1,2-Dichloropropane	1.0	U	
10061-01-5	cis-1,3-Dichloropropene	1.0	บ	
10061-02-6	trans-1,3-Dichloropropene	1.0	U	
75-09-2	Methylene chloride	5.0	ប	
79-34-5	1,1,2,2-Tetrachlorethane	1.0	บ	
127-18-4	Tetrachloroethene	1.0	บ	
71-55-6	1,1,1-Trichloroethane	1.0	U	
79-00-5	1,1,2-Trichloroethane	1.0	ប	
79-01-6	Trichloroethene	1.0	บ	
75-69-4	Trichlorofluoromethane	1.0	บ	
75-01-4	Vinyl chloride	1.0	1.4	

Jul

SURROGATE-Fluorobenzene (QC Limits - 61-133%)

# REPORT OF ANALYTICAL RESULTS PURGEABLE HALOCARBONS

Date Collected: 11/25/96
Date Extracted: N/A

Date Analyzed: 12/02/96 Matrix: Water

Method: EPA 601 (Mod) % Moisture: 100

Sample Group: MC484
Lab Sample ID: MC484004
Lab File 1 ID: N02V011
Lab File 2 ID: N02W011
Dilution Factor: 1.0

Reporting Units: ug/L

CAS NUMBER	COMPOUND NAME	REPORTING LIMIT	RESULT
75-27-4	Bromodichloromethane	1.0	U
75-25-2	Bromoform	1.0	ប
74-83-9	Bromomethane	1.0	ប
56-23-5	Carbon tetrachloride	1.0	ប
108-90-7	Chlorobenzene	1.0	ប
75-00-3	Chloroethane	1.0	ប
110-75-8	2-Chloroethyl vinyl ether	1.0	υ
67-66-3	Chloroform	1.0	ប
74-87-3	Chloromethane	1.0	. υ
124-48-1	Dibromochloromethane	1.0	υ
95-50-1	1,2-Dichlorobenzene	1.0	ប
541-73-1	1,3-Dichlorobenzene	1.0	ប
106-46-7	1,4-Dichlorobenzene	1.0	ប
75-71-8	Dichlorodifluoromethane	1.0	υ
75-34-3	1,1-Dichloroethane	1.0	ប
107-06-2	1,2-Dichloroethane	1.0	υ
75-35-4	1,1-Dichloroethene	1.0	ប
156-59-2	cis-1,2-Dichloroethene	1.0	υ
156-60-5	trans-1,2-Dichloroethene	1.0	U
78-87-5	1,2-Dichloropropane	1.0	ប
10061-01-5	cis-1,3-Dichloropropene	1.0	υ
10061-02-6	trans-1,3-Dichloropropene	1.0	ប
75-09-2	Methylene chloride	5.0	U
79-34-5	1,1,2,2-Tetrachlorethane	1.0	ប
127-18-4	Tetrachloroethene	1.0	ប
71-55-6	1,1,1-Trichloroethane	1.0	ប
79-00-5	1,1,2-Trichloroethane	1.0	U
79-01-6	Trichloroethene	1.0	U
75-69-4	Trichlorofluoromethane	1.0	U
75-01-4	Vinyl chloride	1.0	U

SURROGATE-Fluorobenzene (QC Limits - 61-133%)

65 % Rec.



GC PURGEABLE AROMATICS

## CASE NARRATIVE GC PURGEABLE AROMATICS

QAL Lab	Reference	No./SDG	: MC484		 
Project:	Brown	& Root	Coastal Systems	Station	

#### I. RECEIPT

No exceptions were encountered unless a Sample Receipt Exception Report is attached to the Chain-of-Custody included with this data package.

#### II. HOLDING TIMES

- A. Sample Preparation: N/A
- B. Sample Analysis: All holding times were met.

#### III. METHOD

Preparation:

Cleanup: N/A

Analysis: EPA 602 (Mod)

#### IV. PREPARATION

Not applicable.

#### V. ANALYSIS

- A. Calibration: All acceptance criteria were met.
- B. Blanks: All acceptance criteria were met.
- C. Surrogates: All acceptance criteria were met.
- D. Spikes: All acceptance criteria were met.
- E. Samples: Sample analysis proceeded normally. Primary analysis was performed using a J&W DB-VRX analytical column (75 m x 0.45 mm ID). Confirmational analyses were performed using a Restek Rtx-502.2 analytical column (105 m x 0.53 mm ID).

I certify that this data package is in compliance with the terms and conditions agreed to by the client and QAL, Inc., both technically and for completeness except for the conditions noted above. Release of the data contained in this hardcopy data package has been authorized by the Laboratory Manager or designated person, as verified by the following signature.

SIGNED: Well W. Smiles	DATE:	12-20-96	
Herb Kelly Organic Division Manager			

GC PURGEABLE AROMATICS
Lab Reference No./SDG: MC484

Page 2

#### CASE NARRATIVE Addendum

Sample Information

LAB SAMPLE ID	CLIENT SAMPLE ID	SAMPLE MATRIX	DATE SAMPLED	DATE EXTRACTED	DATE ANALYZED	SAMPLE PH <sup>1</sup>
MC484002	TW01-001B	WATER	11/25/96	N/A	12/02/96	<2
MC484003	TW01-001	WATER	11/25/96	N/A	12/02/96	<2
MC484003MS	TW01-001MS	WATER	11/25/96	N/A	12/02/96	<2
MC484003MSD	TW01-001MSD	WATER	11/25/96	N/A	12/02/96	<2
MC484004	MW04-001	WATER	11/25/96	N/A	12/02/96	<2
X12026B1	VBLK001	WATER	N/A	N/A	12/02/96	N/A

 $<sup>^{\</sup>scriptsize 1}$  Applies to samples designated for purgeable VOA analysis only.

#### REPORT OF ANALYTICAL RESULTS PURGEABLE AROMATICS

Date Collected: 11/25/96 Date Extracted: N/A

Date Analyzed: 12/02/96 Matrix: Water

Method: EPA 602 (Mod)

% Moisture: 100

Sample Group: MC484

Lab Sample ID: MC484002 Lab File 1 ID: N02V012

Lab File 2 ID: N02W012 Dilution Factor: 1.0

Reporting Units: ug/L

CAS NUMBER	COMPOUND NAME	REPORTING LIMIT	RESULT
71-43-2	Benzene	1.0	U
108-88-3	Toluene	1.0	ប
100-41-4	Ethylbenzene	1.0	ט
1330-20-7	Total Xylenes	1.0	บ
N/A	Total Volatile Organic Aromatics	1.0	υ
1634-04-4	Methyl tert-butyl ether	1.0	1.3
	SURROGATE-Fluorobenzene (QC Limits	s - 61-133%)	101 % Rec.



# REPORT OF ANALYTICAL RESULTS PURGEABLE AROMATICS

Date Collected: 11/25/96
Date Extracted: N/A
Date Analyzed: 12/02/96

Matrix: Water

Method: EPA 602 (Mod) % Moisture: 100

Sample Group: MC484 Lab Sample ID: MC484003 Lab File 1 ID: N02V013

Lab File 2 ID: NO2W013
Dilution Factor: 1.0

Reporting Units: ug/L

CAS NUMBER	COMPOUND NAME	REPORTING LIMIT	RESULT
71-43-2	Benzene	1.0	2.2
108-88-3	Toluene	1.0	U
100-41-4	Ethylbenzene	1.0	ប
1330-20-7	Total Xylenes	1.0	ט
N/A	Total Volatile Organic Aromatics	1.0	2.2
1634-04-4	Methyl tert-butyl ether	1.0	U



# REPORT OF ANALYTICAL RESULTS PURGEABLE AROMATICS

Date Collected: 11/25/96

Date Extracted: N/A

Date Analyzed: 12/02/96

Matrix: Water

Method: EPA 602 (Mod)

Sample Group: MC484

Lab Sample ID: MC484004

Lab File 1 ID: N02V011

Lab File 2 ID: N02W011

Dilution Factor: 1.0

% Moisture: 100 Reporting Units: ug/L

CAS NUMBER	COMPOUND NAME	REPORTING LIMIT	RESULT
71-43-2	Benzene	1.0	1.8
108-88-3	Toluene	1.0	U
100-41-4	Ethylbenzene	1.0	4.7
1330-20-7	Total Xylenes	1.0	2.6
N/A	Total Volatile Organic Aromatics	1.0	9.1
1634-04-4	Methyl tert-butyl ether	1.0	T



GC/MS SEMIVOLATILE ORGANICS

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# CASE NARRATIVE GC/MS SEMIVOLATILE ORGANICS

QAL Lab R	eference No./	SDG	MC484	
Project:	Brown &	Root Co	astal Systems	s Station

#### I. RECEIPT

No exceptions were encountered unless a Sample Receipt Exception Report is attached to the Chain-of-Custody included with this data package.

#### II. HOLDING TIMES

- A. Sample Preparation: All holding times were met.
- B. Sample Analysis: All holding times were met.

#### III. METHOD

Preparation: SW-846 3520A

Cleanup: N/A

Analysis: SW-846 8270A

#### IV. PREPARATION

Sample preparation proceeded normally.

#### V. ANALYSIS

- A. Calibration: All acceptance criteria were met.
- B. Blanks: All acceptance criteria were met.
- C. Surrogates: The surrogate recovery for Terphenyl-d14 was lower than QC criteria in samples MC484003MS and MC484003MSD. All other surrogate recoveries were within QC criteria.
- D. Spikes: As requested, the matrix spikes were performed using a sample from sample delivery group MC484 (MC484003MS and MC484003MSD). All acceptance criteria were met. A copy of the results has been included for your review.
- E. Samples: Sample MC484004 was analyzed at a dilution due to nontarget compounds. One set of data is reported for this sample.
- F. Other: Please note that the Form 1's reflect the specified target list.

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GC/MS SEMIVOL	ATILE ORGANICS	
Lab Reference	No./SDG:	MC484
Page 2		

I certify that this data package is in compliance with the terms and conditions agreed to by the client and QAL, Inc., both technically and for completeness except for the conditions noted above. Release of the data contained in this hardcopy data package has been authorized by the Laboratory Manager or designated person, as verified by the following signature.

SIGNED:

DATE: 12-26-96

Ward Dickens

Laboratory Director

GC/MS SEMIVOI	ATILE ORGAN	ICS
Lab Reference	No./SDG:	MC484
Page 3		

# CASE NARRATIVE Addendum

Sample Information

LAB SAMPLE ID	CLIENT SAMPLE ID	SAMPLE MATRIX	DATE SAMPLED	DATE EXTRACTED	DATE ANALYZED	SAMPLE pH <sup>1</sup>
MC484002	TW01-001B	WATER	11/25/96	11/30/96	12/05/96	N/A
MC484003	TW01-001	WATER	11/25/96	11/30/96	12/05/96	N/A
MC484003MS	TW01-001MS	WATER	11/25/96	11/30/96	12/05/96	N/A
MC484003MSD	TW01-001MSD	WATER	11/25/96	11/30/96	12/05/96	N/A
MC484004	MW04-001	WATER	11/25/96	11/30/96	12/05/96	N/A
C11306B2	SBLKRE	WATER	N/A	11/30/96	12/05/96	N/A

<sup>&</sup>lt;sup>1</sup> Applies to samples designated for purgeable VOA analysis only.

EPA SAMPLE NO.

TW01-001B

Lab Name: CH2M HILL Contract: MC484

Lab Code: MGM Case No.: MC484 SAS No.: SDG No.: MC484

Matrix: (soil/water) WATER Lab Sample ID: MC484002

Sample wt/vol: 1000 (q/mL) ML Lab File ID: 05DEC0501005.D

Level: (low/med) LOW Date Received: 11/27/96

% Moisture: not dec. dec. Date Extracted:11/30/96

Extraction: (SepF/Cont/Sonc) CONT Date Analyzed: 12/05/96

GPC Cleanup: (Y/N) N pH: 6.0 Dilution Factor: 1.0

CONCENTRATION UNITS:

CAS NO. COMPOUND (ug/L or ug/Kg) UG/L Q

62-75-9N-Nitrosodimethylamine	10	ט
108-95-2Phenol	10	ן ט
111-44-4bis(2-Chloroethyl)ether	10	ט ט
95-57-82-Chlorophenol	10	Ū
541-73-11,3-Dichlorobenzene	10	ָ װ
106-46-71,4-Dichlorobenzene	10	, ŭ
95-50-11,2-Dichlorobenzene	10	Ŭ
108-60-12,2'-Oxybis(1-chloroprop_(1)	10	Ū
621-64-7Nitroso-di-n-propylamine_	10	ן ט
67-72-1Hexachloroethane	10	ן ט
98-95-3Nitrobenzene	10	ן ט
78-59-1Isophorone	10	ן ט
88-75-52-Nitrophenol	10	الق
105-67-92,4-Dimethylphenol	10	ט .
111-91-1bis (2-Chloroethoxy) methane	10	บี
120-83-22,4-Dichlorophenol	10	บี
120-82-11,2,4-Trichlorobenzene	10	Ü
91-20-3Naphthalene	10	บ
87-68-3Hexachlorobutadiene	10	Ü
59-50-74-Chloro-3-methylphenol	10	ָ ט
88-06-22,4,6-Trichlorophenol	10	Ū.
91-58-72-Chloronaphthalene	10	บ
131-11-3Dimethylphthalate	10	Ü
606-20-22,6-Dinitrotoluene	10	Ü
208-96-8Acenaphthylene	10	บ
83-32-9Acenaphthene	10	Ü
51-28-52,4-Dinitrophenol	50	. Ü
100-02-74-Nitrophenol	50	Ü
121-14-22,4-Dinitrotoluene	10	Ü
84-66-2Diethylphthalate	10	Ü
86-73-7Fluorene	10	ש
7005-72-34-Chlorophenyl-phenylether	10	<u>ט</u>
7003-72-34-Chitotophienyt-phenytechet		
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(1) 2,2'-oxybis(1-Chloropropane) is known as bis(2-Chloroisopropyl) ether

FORM I SV-1

SW846

4 1

Case No.: MC484 SAS No.:

EPA SAMPLE NO.

TW01-001B

Lab Name: CH2M HILL

Contract: MC484

SDG No.: MC484

Matrix: (soil/water) WATER

Lab Sample ID: MC484002

Sample wt/vol:

Lab Code: MGM

1000 (g/mL) ML

Lab File ID: 05DEC0501005.D

Level: (low/med) LOW

Date Received: 11/27/96

% Moisture: not dec.\_\_\_\_\_ dec.\_\_\_\_

Date Extracted:11/30/96

Extraction: (SepF/Cont/Sonc) CONT

Date Analyzed: 12/05/96

GPC Cleanup: (Y/N) N pH: 6.0

Dilution Factor: 1.0

CONCENTRATION UNITS:

CAS NO.

COMPOUND

(uq/L or ug/Kg) UG/L

0

_		•	
534-52-14,6-Dinitro-2-methyl	phenol	50	U
86-30-6N-Nitrosodiphenylami		10	Ū
122-66-71,2-Diphenylhydrazin		10	
101-55-34-Bromophenyl-phenyl		10	l .
118-74-1Hexachlorobenzene		10	I
87-86-5Pentachlorophenol		50	
85-01-8Phenanthrene		10	
120-12-7Anthracene		10	
84-74-2Di-n-butylphthalate	<del></del>	6	1
206-44-0Fluoranthene		10	
92-87-5Benzidine		10	
129-00-0Pyrene		10	_
85-68-7Butylbenzylphthalate		10	1
56-55-3Benzo (a) anthracene		10	1
91-94-13,3'-Dichlorobenzidi	ne	20	)
218-01-9Chrysene		10	
117-81-7bis (2-Ethylhexyl) pht	halate	52	ı
117-84-0Di-n-octylphthalate	-	10	l
205-99-2Benzo (b) fluoranthene	<del></del>	10	1
207-08-9Benzo(k) fluoranthene		10	·
50-32-8Benzo (a) pyrene	·	10	
193-39-5Indeno(1,2,3-cd)pyre	ne l	10	
53-70-3Dibenz (a, h) anthracen		10	_
191-24-2Benzo(g,h,i)perylene		10	U
131-24-2Belizo(g, ii, 1) per yreile	·	10	0
			i

(3) - Cannot be separated from Diphenylamine

FORM I SV-2

SW846

4 3

EPA SAMPLE NO.

TW01-001B

Lab Name: CH2M HILL

Contract: MC484

Lab Code: MGM Case No.: MC484 SAS No.:

SDG No.: MC484

Matrix: (soil/water) WATER

Lab Sample ID: MC484002

Sample wt/vol: 1000 (g/mL) ML

Lab File ID: 05DEC0501005.D

Level: (low/med) LOW

Date Received: 11/27/96

% Moisture: not dec. dec.

Date Extracted:11/30/96

Date Analyzed: 12/05/96

Extraction: (SepF/Cont/Sonc) CONT

GPC Cleanup: (Y/N) N

pH: 6.0

Dilution Factor: 1.0

Number TICs found: 20

CONCENTRATION UNITS: (uq/L or uq/Kq) UG/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
2. 123-42-2 3. 21460-36-6 4. 5. 111-90-0 6. 7. 8. 822-86-6 9. 617-94-7 10. 11. 12. 10546-70-0 13. 14. 15. 16. 17. 18. 19.	Unknown Ketone 2-Pentanone, 4-hydroxy-4-met 2-Propanol, 1-(2-propenyloxy Unknown Ethanol, 2-(2-ethoxyethoxy)- Unknown Unknown Cyclohexane, 1,2-dichloro-, Benzenemethanol, .alpha.,.al Benzene diisocyanato methyl Unknown Benzamide, N-propyl- Unknown Unknown Unknown Unknown Unknown Unknown Unknown Unknown Unknown Unknown Unknown Phthalate Unknown Unknown	4.277 4.435 5.535 5.614 5.721 6.142	2 5 12 4 5 5 2 4 3 4 2 3 6 3 5 2 6 4 10 360	

EPA SAMPLE NO.

TW01-001

Lab Name: CH2M HILL Contract: MC484

Lab Code: MGM Case No.: MC484 SAS No.: SDG No.: MC484

Matrix: (soil/water) WATER Lab Sample ID: MC484003

Lab File ID: 05DEC0601006.D Sample wt/vol: 1000 (g/mL) ML

Level: (low/med) LOW Date Received: 11/27/96

% Moisture: not dec.\_\_\_\_ dec.\_\_\_ Date Extracted:11/30/96

Extraction: (SepF/Cont/Sonc) CONT Date Analyzed: 12/05/96

GPC Cleanup: (Y/N) N pH: 7.0 Dilution Factor: 1.0

CONCENTRATION UNITS:

(1) 2,2'-oxybis(1-Chloropropane) is known as bis(2-Chloroisopropyl) ether

FORM I SV-1

SW846

EPA SAMPLE NO.

TW01-001

Lab Name: CH2M HILL Contract: MC484

Lab Code: MGM Case No.: MC484 SAS No.: SDG No.: MC484

Matrix: (soil/water) WATER Lab Sample ID: MC484003

Sample wt/vol: 1000 (g/mL) ML Lab File ID: 05DEC0601006.D

Level: (low/med) LOW Date Received: 11/27/96

% Moisture: not dec.\_\_\_\_\_ dec.\_\_\_\_ Date Extracted:11/30/96

Extraction: (SepF/Cont/Sonc) CONT Date Analyzed: 12/05/96

GPC Cleanup: (Y/N) N pH: 7.0 Dilution Factor: 1.0

CONCENTRATION UNITS:

CAS NO.	COMPOUND	(ug/L or ug		Q
86-30-6 122-66-7 101-55-3 118-74-1 87-86-5 85-01-8 120-12-7 84-74-2 206-44-0 92-87-5 129-00-0 85-68-7 56-55-3 91-94-1 218-01-9 117-81-7 117-84-0 205-99-2 207-08-9 50-32-8 193-39-5 53-70-3	4,6-Dinitro-2N-Nitrosodiph1,2-Diphenylh4-Bromophenyl4-BromophenylPentachlorophPhenanthreneAnthraceneDi-n-butylphtFluorantheneBenzidinePyreneButylbenzylphBenzo(a) anthr3,3'-DichloroChrysenebis(2-EthylheDi-n-octylphtBenzo(b) fluorBenzo(a) pyrenIndeno(1,2,3Dibenz(a,h) andBenzo(g,h,i) p	menylamine(3) mydrazinephenyletherphenyletherphenylether	50 10 10 10 10 50 10 10 10 10 10 10 10 10 10 10 10 10	ממממממממממממממממממ
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(3) - Cannot be separated from Diphenylamine

SW846

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EPA SAMPLE NO.

TW01-001

Lab Name: CH2M HILL

Contract: MC484

Lab Code: MGM Case No.: MC484 SAS No.: SDG No.: MC484

Matrix: (soil/water) WATER

Lab Sample ID: MC484003

Sample wt/vol: 1000 (g/mL) ML

Lab File ID: 05DEC0601006.D

Level: (low/med) LOW

Date Received: 11/27/96

% Moisture: not dec.\_\_\_\_ dec.

Date Extracted:11/30/96

Extraction: (SepF/Cont/Sonc) CONT

Date Analyzed: 12/05/96

GPC Cleanup: (Y/N) N

pH: 7.0

Dilution Factor: 1.0

CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L

Number TICs found: 20

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1. 2. 123-42-2 3. 4. 112-36-7 5. 822-86-6 6. 7. 2039-89-6 8. 9. 10. 11. 12. 13. 14. 15. 16. 480-63-7 17. 18. 19. 20. 21. 22. 23. 24. 25. 26. 27. 28. 29. 30.	Unknown Ketone 2-Pentanone, 4-hydroxy-4-met Unknown Ethane, 1,1'-oxybis[2-ethoxy Cyclohexane, 1,2-dichloro-, Unknown Benzene, 2-ethenyl-1,4-dimet Unknown	4.434 5.542 6.142 6.321 6.799 6.942 7.057 7.178 7.614 8.250 8.343 8.400 8.457	6 22 4 4 5 4 3 4 4 3 4 6 7 12	TE NOTE OF THE PROPERTY OF THE

٤.:

EPA SAMPLE NO.

MW04-001

Lab Name: CH2M HILL Contract: MC484

Lab Code: MGM Case No.: MC484 SAS No.: SDG No.: MC484

Matrix: (soil/water) WATER Lab Sample ID: MC484004

Sample wt/vol: 1000 (g/mL) ML Lab File ID: 05DEC0701007.D

Level: (low/med) LOW Date Received: 11/27/96

% Moisture: not dec. dec. Date Extracted:11/30/96

Extraction: (SepF/Cont/Sonc) CONT Date Analyzed: 12/05/96

GPC Cleanup: (Y/N) N pH: 7.0 Dilution Factor: 10.0

> CONCENTRATION UNITS: CAS NO. COMPOUND (ug/L or ug/Kg) UG/L 0

CAS NO.	COMPOUND (ag/L of ag	J/ Kg/ OG/ L	Q
62-75-9	N-Nitrosodimethylamine	100	ט
108-95-2		100	ע
	bis(2-Chloroethyl)ether	100	ע
	2-Chlorophenol	100	ן ט
	1,3-Dichlorobenzene	100	ן ט
	1,4-Dichlorobenzene	100	U
	1,2-Dichlorobenzene	100	ਂ ਹ
	2,2'-Oxybis(1-chloroprop_(1)	100	Ū
621-64-7	N-Nitroso-di-n-propylamine	100	U
	Hexachloroethane	100	U
	Nitrobenzene	100	ט
78-59-1	Isophorone	100	ט
	2-Nitrophenol	100	U
105-67-9	2,4-Dimethylphenol	100	Ū
111-91-1	bis(2-Chloroethoxy)methane	100	U
120-83-2	2,4-Dichlorophenol	100	Ū
	1,2,4-Trichlorobenzene	100	J
91-20-3	Naphthalene	55	J
87-68-3	Hexachlorobutadiene	100	U
59-50-7	4-Chloro-3-methylphenol	100	U
	2,4,6-Trichlorophenol	100	U
91-58-7	2-Chloronaphthalene	100	U
	Dimethylphthalate	100	U
	2,6-Dinitrotoluene	100	บ
	Acenaphthylene	100	U
83-32-9	Acenaphthene	53	J
	2,4-Dinitrophenol	500	U
	4-Nitrophenol	500	Ū
	2,4-Dinitrotoluene	100	ַ ט
	Diethylphthalate	100	U
	Fluorene	_  81	J
7005-72-3	4-Chlorophenyl-phenylether_	100	띠
		_	ll

(1) 2,2'-oxybis(1-Chloropropane) is known as bis(2-Chloroisopropyl) ether

FORM I SV-1

SW846